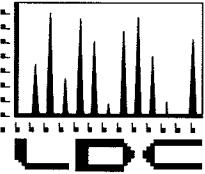


# **Appendix E**

## **Data Validation Reports**

**CDM**



**LABORATORY DATA CONSULTANTS, INC.**  
7750 El Camino Real, Suite 2L Carlsbad, CA 92009 Phone: 760/634-0437 Fax: 760/634-0439

Tait Environmental Management, Inc.  
701 N. Park Center Drive  
Santa Ana, CA 92705  
ATTN: Mr. Matt Hillman

May 16, 2005

SUBJECT: Boeing Realty Corp., Bldg. C-6 Facility, Data Validation

Dear Mr. Hillman,

Enclosed are the final validation reports for the fractions listed below. These SDGs were received on October 20, 2005. Attachment 1 is a summary of the samples that were reviewed for each analysis.

**LDC Project # 14879:**

<b><u>SDG #</u></b>	<b><u>Fraction</u></b>
IPC2080, IPC2190,	Volatiles, Manganese, Wet Chemistry, Fixed
IPC2325	Gases

The data validation was performed under Tier 1, Tier 2 and Tier 3 guidelines. The analyses were validated using the following documents, as applicable to each method:

- USEPA, Contract Laboratory Program National Functional Guidelines for Organic Data Review, October 1999
- USEPA, Contract Laboratory Program National Functional Guidelines for Inorganic Data Review, October 2004
- EPA SW 846, Third Edition, Test Methods for Evaluating Solid Waste, update 1, July 1992; update IIA, August 1993; update II, September 1994; update IIB, January 1995; update III, December 1996; update IIIA, April 1998

Please feel free to contact us if you have any questions.

Sincerely,

*E. Pautz Jr.*  
Stella S. Cuenco  
Project Manager/Senior Chemist

LDC #14879 (Tait Environmental Management, Inc. / Boeing Realty Corp., Former C-6 Facility, Torrance, CA)

Shaded cells indicate Tier III validation (all other cells are Tier II validation). Sample counts do not include MS, MSD, or DLIP's

**Boeing Realty Corp., Former C-6 Facility, Torrance CA**  
**Data Validation Reports**  
**LDC# 14879**

**Volatiles**

E D C

**Laboratory Data Consultants, Inc.**  
**Data Validation Report**

**Project/Site Name:** Boeing Realty Corp, Former C-6 Facility, Torrance CA  
**Collection Date:** March 20, 2006  
**LDC Report Date:** May 15, 2006  
**Matrix:** Water  
**Parameters:** Volatiles  
**Validation Level:** Tier 1  
**Laboratory:** Del Mar Analytical  
**Sample Delivery Group (SDG):** IPC2080

**Sample Identification**

MWB003\_WG032006\_0001  
MWC011\_WG032006\_0001  
IRZMW002B\_WG032006\_0001  
IRZMW002B\_WG032006\_0002  
IRZB0081\_WG032006\_0001  
MWB019\_WG032006\_0001  
MWB019\_WG032006\_0001MS  
MWB019\_WG032006\_0001MSD  
IRZMW002B\_WG032006\_0002MS  
IRZMW002B\_WG032006\_0002MSD

## **Introduction**

This data review covers 10 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 8260B for Volatiles.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review (October 1999) as there are no current guidelines for the method stated above.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified a P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blank results are summarized in Section V.

Field duplicates are summarized in Section XVI.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- U     Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- J     Indicates an estimated value.
- R     Quality control indicates the data is not usable.
- N     Presumptive evidence of presence of the constituent.
- UJ    Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A     Indicates the finding is based upon technical validation criteria.
- P     Indicates the finding is related to a protocol/contractual deviation.
- None    Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodices were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. GC/MS Instrument Performance Check**

Instrument performance check data were not reviewed for Tier 1.

## **III. Initial Calibration**

Initial calibration data were not reviewed for Tier 1.

## **IV. Continuing Calibration**

Continuing calibration data were not reviewed for Tier 1.

## **V. Blanks**

Method blanks were reviewed for each matrix as applicable. No volatile contaminants were found in the method blanks with the following exceptions:

Method Blank ID	Analysis Date	Compound TIC (RT in minutes)	Concentration	Associated Samples
6C29002-BLK	3/29/06	Tetrahydrofuran	3.51 ug/L	MWB003_WG032006_0001 MWC011_WG032006_0001 IRZMW002B_WG032006_0001 IRZMW002B_WG032006_0002 IRZB0081_WG032006_0001
6C30003-BLK1	3/30/06	Tetrahydrofuran	7.19 ug/L	MWB019_WG032006_0001

Sample concentrations were compared to concentrations detected in the method blanks. The sample concentrations were either not detected or were significantly greater (>10X for common contaminants, >5X for other contaminants) than the concentrations found in the associated method blanks.

## **VI. Surrogate Spikes**

Surrogates were added to all samples and blanks as required by the method. All surrogate recoveries (%R) were within QC limits with the following exceptions:

Sample	Surrogate	%R (Limits)	Compound	Flag	A or P
MWC011_WG032006_0001	Dibromofluoromethane	122 (80-120)	Tetrahydrofuran Iodomethane	J (all detects) UJ (all non-detects) J (all detects) UJ (all non-detects)	A
IRZB0081_WG032006_0001	Dibromofluoromethane	121 (80-120)	Tetrahydrofuran Iodomethane	J (all detects) UJ (all non-detects) J (all detects) UJ (all non-detects)	A

## VII. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits with the following exceptions:

Spike ID (Associated Samples)	Compound	MS (%R) (Limits)	MSD (%R) (Limits)	RPD (Limits)	Flag	A or P
MWB019_WG032006_0001MS/MSD (MWB019_WG032006_0001)	n-Butylbenzene 1,2-Dichloropropane Ethylbenzene Hexachlorobutadiene	140 (65-135) 130 (60-125) 132 (65-130) 136 (60-135)	- - - -	- - - -	J (all detects) J (all detects) J (all detects) J (all detects)	A
MWB019_WG032006_0001MS/MSD (MWB019_WG032006_0001)	Styrene 1,2,4-Trimethylbenzene	- -	23 (45-145) 52 (55-130)	75 ( $\leq$ 30) 46 ( $\leq$ 25)	J (all detects) UJ (all non-detects) J (all detects) UJ (all non-detects)	A
IRZMW002B_WG032006_0002MS/MSD (IRZMW002B_WG032006_0002RE2)	Acetone	-	-	42 ( $\leq$ 35)	J (all detects) UJ (all non-detects)	A

## VIII. Laboratory Control Samples (LCS)

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits with the following exceptions:

LCS ID	Compound	%R (Limits)	Associated Samples	Flag	A or P
6C30003-BS1	4-Methyl-2-pentanone	136 (40-135)	MWB019_WG032006_0001 6C30003-BLK1	J (all detects)	P

## IX. Regional Quality Assurance and Quality Control

Not applicable.

## **X. Internal Standards**

Internal standards data were not reviewed for Tier 1.

## **XI. Target Compound Identifications**

Raw data were not reviewed for this SDG.

## **XII. Compound Quantitation and CRQLs**

Raw data were not reviewed for this SDG.

## **XIII. Tentatively Identified Compounds (TICs)**

Raw data were not reviewed for this SDG.

## **XIV. System Performance**

Raw data were not reviewed for this SDG.

## **XV. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **XVI. Field Duplicates**

Samples IRZMW002B\_WG032006\_0001 and IRZMW002B\_WG032006\_0002 were identified as field duplicates. No volatiles were detected in any of the samples with the following exceptions:

Compound	Concentration (ug/L)		RPD
	IRZMW002B_WG032006_0001	IRZMW002B_WG032006_0002	
Trichloroethene	300	330	10
Chlorobenzene	0.42	0.48	13
Chloroform	1.5	1.3	14
1,1-Dichloroethene	6.2	3.9	46
cis-1,2-Dichloroethene	110	94	16
Toluene	1.4	1.3	7

Compound	Concentration (ug/L)		RPD
	IRZMW002B_WG032006_0001	IRZMW002B_WG032006_0002	
trans-1,2-Dichloroethene	1.0U	4.2	200

## XVII. Field Blanks

No field blanks were identified in this SDG.

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Volatiles - Data Qualification Summary - SDG IPC2080**

SDG	Sample	Compound	Flag	A or P	Reason
IPC2080	MWC011_WG032006_0001 IRZB0081_WG032006_0001	Tetrahydrofuran Iodomethane	J (all detects) UJ (all non-detects) J (all detects) UJ (all non-detects)	A	Surrogate recovery (%R)
	MWB019_WG032006_0001	n-Butylbenzene 1,2-Dichloropropane Ethylbenzene Hexachlorobutadiene	J (all detects) J (all detects) J (all detects) J (all detects)	A	Matrix spike/Matrix spike duplicates (%R)
IPC2080	MWB019_WG032006_0001	Styrene 1,2,4-Trimethylbenzene	J (all detects) UJ (all non-detects) J (all detects) UJ (all non-detects)	A	Matrix spike/Matrix spike duplicates (%R)(RPD)
IPC2080	IRZMW002B_WG032006_0002	Acetone	J (all detects) UJ (all non-detects)	A	Matrix spike/Matrix spike duplicates (RPD)
IPC2080	MWB019_WG032006_0001	4-Methyl-2-pentanone	J (all detects)	P	Laboratory control samples (%R)

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Volatiles - Laboratory Blank Data Qualification Summary - SDG IPC2080**

No Sample Data Qualified in this SDG



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 2520 E. Sunset Rd. #3, Las Vegas, NV 89120 (702) 798-3620 FAX (702) 798-3621

TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2080

Sampled: 03/20/06  
 Received: 03/20/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2080-06 (MWB003_WG032006_0001 - Water)</b>									
<b>Reporting Units: ug/l</b>									
Acetone	EPA 8260B	6C29002	4.5	10	ND	1	03/29/06	03/29/06	
Benzene	EPA 8260B	6C29002	0.28	1.0	9.3	1	03/29/06	03/29/06	
Bromobenzene	EPA 8260B	6C29002	0.27	1.0	ND	1	03/29/06	03/29/06	
Bromoform	EPA 8260B	6C29002	0.32	1.0	ND	1	03/29/06	03/29/06	
Bromomethane	EPA 8260B	6C29002	0.30	1.0	ND	1	03/29/06	03/29/06	
2-Butanone (MEK)	EPA 8260B	6C29002	3.8	5.0	ND	1	03/29/06	03/29/06	
n-Butylbenzene	EPA 8260B	6C29002	0.37	1.0	ND	1	03/29/06	03/29/06	
sec-Butylbenzene	EPA 8260B	6C29002	0.25	1.0	ND	1	03/29/06	03/29/06	
tert-Butylbenzene	EPA 8260B	6C29002	0.22	1.0	ND	1	03/29/06	03/29/06	
Carbon Disulfide	EPA 8260B	6C29002	0.48	1.0	ND	1	03/29/06	03/29/06	
Carbon tetrachloride	EPA 8260B	6C29002	0.28	0.50	ND	1	03/29/06	03/29/06	
Chlorobenzene	EPA 8260B	6C29002	0.36	1.0	ND	1	03/29/06	03/29/06	
Chloroethane	EPA 8260B	6C29002	0.40	2.0	ND	1	03/29/06	03/29/06	
<b>Chloroform</b>	EPA 8260B	6C29002	0.33	1.0	25	1	03/29/06	03/29/06	
Chloromethane	EPA 8260B	6C29002	0.30	2.0	ND	1	03/29/06	03/29/06	
2-Chlorotoluene	EPA 8260B	6C29002	0.28	1.0	ND	1	03/29/06	03/29/06	
4-Chlorotoluene	EPA 8260B	6C29002	0.29	1.0	ND	1	03/29/06	03/29/06	
1,2-Dibromo-3-chloropropane	EPA 8260B	6C29002	0.92	2.0	ND	1	03/29/06	03/29/06	
Dibromochloromethane	EPA 8260B	6C29002	0.28	1.0	ND	1	03/29/06	03/29/06	
1,2-Dibromoethane (EDB)	EPA 8260B	6C29002	0.32	1.0	ND	1	03/29/06	03/29/06	
1,4-Dichlorobenzene	EPA 8260B	6C29002	0.37	1.0	ND	1	03/29/06	03/29/06	
1,2-Dichlorobenzene	EPA 8260B	6C29002	0.32	1.0	ND	1	03/29/06	03/29/06	
1,3-Dichlorobenzene	EPA 8260B	6C29002	0.35	1.0	ND	1	03/29/06	03/29/06	
Dichlorodifluoromethane	EPA 8260B	6C29002	0.79	1.0	ND	1	03/29/06	03/29/06	
<b>1,2-Dichloroethane</b>	EPA 8260B	6C29002	0.28	0.50	41	1	03/29/06	03/29/06	
<b>1,1-Dichloroethane</b>	EPA 8260B	6C29002	0.27	1.0	89	1	03/29/06	03/29/06	
<b>cis-1,2-Dichloroethene</b>	EPA 8260B	6C29002	0.32	1.0	130	1	03/29/06	03/29/06	
<b>trans-1,2-Dichloroethene</b>	EPA 8260B	6C29002	0.27	1.0	98	1	03/29/06	03/29/06	
1,2-Dichloropropane	EPA 8260B	6C29002	0.35	1.0	ND	1	03/29/06	03/29/06	
2,2-Dichloropropane	EPA 8260B	6C29002	0.34	1.0	ND	1	03/29/06	03/29/06	
cis-1,3-Dichloropropene	EPA 8260B	6C29002	0.22	0.50	ND	1	03/29/06	03/29/06	
1,1-Dichloropropene	EPA 8260B	6C29002	0.28	1.0	ND	1	03/29/06	03/29/06	
trans-1,3-Dichloropropene	EPA 8260B	6C29002	0.32	0.50	ND	1	03/29/06	03/29/06	
Ethylbenzene	EPA 8260B	6C29002	0.25	1.0	ND	1	03/29/06	03/29/06	
Hexachlorobutadiene	EPA 8260B	6C29002	0.38	1.0	ND	1	03/29/06	03/29/06	
2-Hexanone	EPA 8260B	6C29002	2.6	6.0	ND	1	03/29/06	03/29/06	
Iodomethane	EPA 8260B	6C29002	1.0	2.0	ND	1	03/29/06	03/29/06	
Isopropylbenzene	EPA 8260B	6C29002	0.25	1.0	ND	1	03/29/06	03/29/06	
p-Isopropyltoluene	EPA 8260B	6C29002	0.28	1.0	ND	1	03/29/06	03/29/06	

**Del Mar Analytical - Irvine**  
 Michele Chamberlin  
 Project Manager

KOTRUC

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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2080

Sampled: 03/20/06  
 Received: 03/20/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2080-06 (MWB003_WG032006_0001 - Water) - cont.</b>									
Reporting Units: ug/l									
Methyl-tert-butyl Ether (MTBE)	EPA 8260B	6C29002	0.32	1.0	ND	1	03/29/06	03/29/06	
<b>Methylene chloride</b>	EPA 8260B	6C29002	0.70	1.0	<b>2.2</b>	1	03/29/06	03/29/06	
4-Methyl-2-pentanone (MIBK)	EPA 8260B	6C29002	3.5	5.0	ND	1	03/29/06	03/29/06	
n-Propylbenzene	EPA 8260B	6C29002	0.27	1.0	ND	1	03/29/06	03/29/06	
Styrene	EPA 8260B	6C29002	0.16	1.0	ND	1	03/29/06	03/29/06	
1,1,1,2-Tetrachloroethane	EPA 8260B	6C29002	0.27	1.0	ND	1	03/29/06	03/29/06	
1,1,2,2-Tetrachloroethane	EPA 8260B	6C29002	0.24	1.0	ND	1	03/29/06	03/29/06	
<b>Tetrachloroethene</b>	EPA 8260B	6C29002	0.32	1.0	<b>0.50</b>	1	03/29/06	03/29/06	J
Tetrahydrofuran (THF)	EPA 8260B	6C29002	2.3	10	ND	1	03/29/06	03/29/06	
Toluene	EPA 8260B	6C29002	0.36	1.0	ND	1	03/29/06	03/29/06	
1,2,3-Trichlorobenzene	EPA 8260B	6C29002	0.45	1.0	ND	1	03/29/06	03/29/06	
1,2,4-Trichlorobenzene	EPA 8260B	6C29002	0.48	1.0	ND	1	03/29/06	03/29/06	
<b>1,1,2-Trichloroethane</b>	EPA 8260B	6C29002	0.30	1.0	<b>27</b>	1	03/29/06	03/29/06	
<b>1,1,1-Trichloroethane</b>	EPA 8260B	6C29002	0.30	1.0	<b>1.9</b>	1	03/29/06	03/29/06	
Trichlorofluoromethane	EPA 8260B	6C29002	0.34	2.0	ND	1	03/29/06	03/29/06	
1,2,3-Trichloropropane	EPA 8260B	6C29002	0.40	1.0	ND	1	03/29/06	03/29/06	
1,2,4-Trimethylbenzene	EPA 8260B	6C29002	0.23	1.0	ND	1	03/29/06	03/29/06	
1,3,5-Trimethylbenzene	EPA 8260B	6C29002	0.26	1.0	ND	1	03/29/06	03/29/06	
Vinyl acetate	EPA 8260B	6C29002	1.7	6.0	ND	1	03/29/06	03/29/06	
Vinyl chloride	EPA 8260B	6C29002	0.26	0.50	ND	1	03/29/06	03/29/06	
Xylenes, Total	EPA 8260B	6C29002	0.90	1.0	ND	1	03/29/06	03/29/06	
<i>Surrogate: 4-Bromofluorobenzene (80-120%)</i>					96 %				
<i>Surrogate: Dibromofluoromethane (80-120%)</i>					116 %				
<i>Surrogate: Toluene-d8 (80-120%)</i>					103 %				

**Del Mar Analytical - Irvine**  
 Michele Chamberlin  
 Project Manager

7/25/06

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TAIT Environmental/Boeing  
701 N. Parkcenter Drive  
Santa Ana, CA 92705  
Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
EM2303D/PHASE 01  
Report Number: IPC2080

Sampled: 03/20/06  
Received: 03/20/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2080-06RE1 (MWB003_WG032006_0001 - Water) - cont.</b>									
Reporting Units: ug/l									
1,1-Dichloroethene	EPA 8260B	6C31009	42	100	4800	100	03/31/06	03/31/06	
Trichloroethene	EPA 8260B	6C31009	26	100	4100	100	03/31/06	03/31/06	
Surrogate: 4-Bromofluorobenzene (80-120%)									
Surrogate: Dibromofluoromethane (80-120%)									
Surrogate: Toluene-d8 (80-120%)									

Del Mar Analytical - Irvine  
Michele Chamberlin  
Project Manager

*K. Chamberlin*

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BOE-C6-0050665



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 2520 E. Sunset Rd. #3, Las Vegas, NV 89120 (702) 798-3620 FAX (702) 798-3621

TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
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 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2080

Sampled: 03/20/06  
 Received: 03/20/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2080-07 (MWC011_WG032006_0001 - Water)</b>									
Reporting Units: ug/l									
Iodomethane	EPA 8260B	6C29002	1.0	2.0	ND	1	03/29/06	03/29/06	
Tetrahydrofuran (THF)	EPA 8260B	6C29002	2.3	10	ND	1	03/29/06	03/29/06	
<i>Surrogate: 4-Bromofluorobenzene (80-120%)</i>									
<i>Surrogate: Dibromofluoromethane (80-120%)</i>									
<i>Surrogate: Toluene-d8 (80-120%)</i>									
<b>Sample ID: IPC2080-07RE1 (MWC011_WG032006_0001 - Water)</b>									
Reporting Units: ug/l									
Acetone	EPA 8260B	6C31009	4.5	10	ND	1	03/31/06	03/31/06	
<b>Benzene</b>	EPA 8260B	6C31009	0.28	1.0	<b>0.30</b>	1	03/31/06	03/31/06	J
Bromobenzene	EPA 8260B	6C31009	0.27	1.0	ND	1	03/31/06	03/31/06	
Bromoform	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
Bromochloromethane	EPA 8260B	6C31009	0.30	1.0	ND	1	03/31/06	03/31/06	
Bromodichloromethane	EPA 8260B	6C31009	0.30	1.0	ND	1	03/31/06	03/31/06	
Bromoform	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
Bromomethane	EPA 8260B	6C31009	0.42	1.0	ND	1	03/31/06	03/31/06	
2-Butanone (MEK)	EPA 8260B	6C31009	3.8	5.0	ND	1	03/31/06	03/31/06	
n-Butylbenzene	EPA 8260B	6C31009	0.37	1.0	ND	1	03/31/06	03/31/06	
sec-Butylbenzene	EPA 8260B	6C31009	0.25	1.0	ND	1	03/31/06	03/31/06	
tert-Butylbenzene	EPA 8260B	6C31009	0.22	1.0	ND	1	03/31/06	03/31/06	
Carbon Disulfide	EPA 8260B	6C31009	0.48	1.0	ND	1	03/31/06	03/31/06	
Carbon tetrachloride	EPA 8260B	6C31009	0.28	0.50	ND	1	03/31/06	03/31/06	
Chlorobenzene	EPA 8260B	6C31009	0.36	1.0	ND	1	03/31/06	03/31/06	
Chloroethane	EPA 8260B	6C31009	0.40	2.0	ND	1	03/31/06	03/31/06	
Chloroform	EPA 8260B	6C31009	0.33	1.0	ND	1	03/31/06	03/31/06	
Chloromethane	EPA 8260B	6C31009	0.30	2.0	ND	1	03/31/06	03/31/06	
2-Chlorotoluene	EPA 8260B	6C31009	0.28	1.0	ND	1	03/31/06	03/31/06	
4-Chlorotoluene	EPA 8260B	6C31009	0.29	1.0	ND	1	03/31/06	03/31/06	
1,2-Dibromo-3-chloropropane	EPA 8260B	6C31009	0.92	2.0	ND	1	03/31/06	03/31/06	
Dibromochloromethane	EPA 8260B	6C31009	0.28	1.0	ND	1	03/31/06	03/31/06	
1,2-Dibromoethane (EDB)	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
1,4-Dichlorobenzene	EPA 8260B	6C31009	0.37	1.0	ND	1	03/31/06	03/31/06	
1,2-Dichlorobenzene	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
1,3-Dichlorobenzene	EPA 8260B	6C31009	0.35	1.0	ND	1	03/31/06	03/31/06	
Dichlorodifluoromethane	EPA 8260B	6C31009	0.79	1.0	ND	1	03/31/06	03/31/06	
<b>1,2-Dichloroethane</b>	EPA 8260B	6C31009	0.28	0.50	<b>0.60</b>	1	03/31/06	03/31/06	
<b>1,1-Dichloroethane</b>	EPA 8260B	6C31009	0.27	1.0	<b>2.5</b>	1	03/31/06	03/31/06	
<b>1,1-Dichloroethene</b>	EPA 8260B	6C31009	0.42	1.0	<b>59</b>	1	03/31/06	03/31/06	
<b>cis-1,2-Dichloroethene</b>	EPA 8260B	6C31009	0.32	1.0	<b>51</b>	1	03/31/06	03/31/06	
<b>trans-1,2-Dichloroethene</b>	EPA 8260B	6C31009	0.27	1.0	<b>1.4</b>	1	03/31/06	03/31/06	
1,2-Dichloropropane	EPA 8260B	6C31009	0.35	1.0	ND	1	03/31/06	03/31/06	
2,2-Dichloropropane	EPA 8260B	6C31009	0.34	1.0	ND	1	03/31/06	03/31/06	

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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2080

Sampled: 03/20/06  
 Received: 03/20/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2080-07RE1 (MWC011_WG032006_0001 - Water) - cont.</b>									
<b>Reporting Units: ug/l</b>									
cis-1,3-Dichloropropene	EPA 8260B	6C31009	0.22	0.50	ND	1	03/31/06	03/31/06	
1,1-Dichloropropene	EPA 8260B	6C31009	0.28	1.0	ND	1	03/31/06	03/31/06	
trans-1,3-Dichloropropene	EPA 8260B	6C31009	0.32	0.50	ND	1	03/31/06	03/31/06	
Ethylbenzene	EPA 8260B	6C31009	0.25	1.0	ND	1	03/31/06	03/31/06	
Hexachlorobutadiene	EPA 8260B	6C31009	0.38	1.0	ND	1	03/31/06	03/31/06	
2-Hexanone	EPA 8260B	6C31009	2.6	6.0	ND	1	03/31/06	03/31/06	
Isopropylbenzene	EPA 8260B	6C31009	0.25	1.0	ND	1	03/31/06	03/31/06	
p-Isopropyltoluene	EPA 8260B	6C31009	0.28	1.0	ND	1	03/31/06	03/31/06	
Methyl-tert-butyl Ether (MTBE)	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
<b>Methylene chloride</b>	EPA 8260B	6C31009	0.70	1.0	<b>1.5</b>	1	03/31/06	03/31/06	
4-Methyl-2-pentanone (MIBK)	EPA 8260B	6C31009	3.5	5.0	ND	1	03/31/06	03/31/06	
n-Propylbenzene	EPA 8260B	6C31009	0.27	1.0	ND	1	03/31/06	03/31/06	
Styrene	EPA 8260B	6C31009	0.16	1.0	ND	1	03/31/06	03/31/06	
1,1,1,2-Tetrachloroethane	EPA 8260B	6C31009	0.27	1.0	ND	1	03/31/06	03/31/06	
1,1,2,2-Tetrachloroethane	EPA 8260B	6C31009	0.24	1.0	ND	1	03/31/06	03/31/06	
Tetrachloroethene	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
Toluene	EPA 8260B	6C31009	0.36	1.0	ND	1	03/31/06	03/31/06	
1,2,3-Trichlorobenzene	EPA 8260B	6C31009	0.45	1.0	ND	1	03/31/06	03/31/06	
1,2,4-Trichlorobenzene	EPA 8260B	6C31009	0.48	1.0	ND	1	03/31/06	03/31/06	
1,1,2-Trichloroethane	EPA 8260B	6C31009	0.30	1.0	ND	1	03/31/06	03/31/06	
1,1,1-Trichloroethane	EPA 8260B	6C31009	0.30	1.0	ND	1	03/31/06	03/31/06	
<b>Trichloroethene</b>	EPA 8260B	6C31009	0.26	1.0	<b>33</b>	1	03/31/06	03/31/06	
Trichlorofluoromethane	EPA 8260B	6C31009	0.34	2.0	ND	1	03/31/06	03/31/06	
1,2,3-Trichloropropane	EPA 8260B	6C31009	0.40	1.0	ND	1	03/31/06	03/31/06	
1,2,4-Trimethylbenzene	EPA 8260B	6C31009	0.23	1.0	ND	1	03/31/06	03/31/06	
1,3,5-Trimethylbenzene	EPA 8260B	6C31009	0.26	1.0	ND	1	03/31/06	03/31/06	
Vinyl acetate	EPA 8260B	6C31009	1.7	6.0	ND	1	03/31/06	03/31/06	
Vinyl chloride	EPA 8260B	6C31009	0.26	0.50	ND	1	03/31/06	03/31/06	
Xylenes, Total	EPA 8260B	6C31009	0.90	1.0	ND	1	03/31/06	03/31/06	
Surrogate: 4-Bromofluorobenzene (80-120%)									
103 %									
Surrogate: Dibromofluoromethane (80-120%)									
107 %									
Surrogate: Toluene-d8 (80-120%)									
107 %									

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 Project Manager

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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2080

Sampled: 03/20/06  
 Received: 03/20/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2080-08 (IRZMW002B_WG032006_0001 - Water)</b>									
<b>Reporting Units: ug/l</b>									
Iodomethane	EPA 8260B	6C29002	1.0	2.0	ND	1	03/29/06	03/29/06	
Tetrahydrofuran (THF)	EPA 8260B	6C29002	2.3	10	ND	1	03/29/06	03/29/06	
<i>Surrogate: 4-Bromofluorobenzene (80-120%)</i>					95 %				
<i>Surrogate: Dibromofluoromethane (80-120%)</i>					120 %				
<i>Surrogate: Toluene-d8 (80-120%)</i>					102 %				
<b>Sample ID: IPC2080-08RE1 (IRZMW002B_WG032006_0001 - Water)</b>									
<b>Reporting Units: ug/l</b>									
Trichloroethene	EPA 8260B	6C31009	1.3	5.0	300	5	03/31/06	03/31/06	
<i>Surrogate: 4-Bromofluorobenzene (80-120%)</i>					110 %				
<i>Surrogate: Dibromofluoromethane (80-120%)</i>					110 %				
<i>Surrogate: Toluene-d8 (80-120%)</i>					108 %				
<b>Sample ID: IPC2080-08RE2 (IRZMW002B_WG032006_0001 - Water)</b>									
<b>Reporting Units: ug/l</b>									
Acetone	EPA 8260B	6C31009	4.5	10	ND	1	03/31/06	03/31/06	
Benzene	EPA 8260B	6C31009	0.28	1.0	ND	1	03/31/06	03/31/06	
Bromobenzene	EPA 8260B	6C31009	0.27	1.0	ND	1	03/31/06	03/31/06	
Bromoform	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
Bromochloromethane	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
Bromodichloromethane	EPA 8260B	6C31009	0.30	1.0	ND	1	03/31/06	03/31/06	
Bromoform	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
Bromomethane	EPA 8260B	6C31009	0.42	1.0	ND	1	03/31/06	03/31/06	
2-Butanone (MEK)	EPA 8260B	6C31009	3.8	5.0	ND	1	03/31/06	03/31/06	
n-Butylbenzene	EPA 8260B	6C31009	0.37	1.0	ND	1	03/31/06	03/31/06	
sec-Butylbenzene	EPA 8260B	6C31009	0.25	1.0	ND	1	03/31/06	03/31/06	
tert-Butylbenzene	EPA 8260B	6C31009	0.22	1.0	ND	1	03/31/06	03/31/06	
Carbon Disulfide	EPA 8260B	6C31009	0.48	1.0	ND	1	03/31/06	03/31/06	
Carbon tetrachloride	EPA 8260B	6C31009	0.28	0.50	ND	1	03/31/06	03/31/06	
<b>Chlorobenzene</b>	EPA 8260B	6C31009	0.36	1.0	0.42	1	03/31/06	03/31/06	J
Chloroethane	EPA 8260B	6C31009	0.40	2.0	ND	1	03/31/06	03/31/06	
<b>Chloroform</b>	EPA 8260B	6C31009	0.33	1.0	1.5	1	03/31/06	03/31/06	
Chloromethane	EPA 8260B	6C31009	0.30	2.0	ND	1	03/31/06	03/31/06	
2-Chlorotoluene	EPA 8260B	6C31009	0.28	1.0	ND	1	03/31/06	03/31/06	
4-Chlorotoluene	EPA 8260B	6C31009	0.29	1.0	ND	1	03/31/06	03/31/06	
1,2-Dibromo-3-chloropropane	EPA 8260B	6C31009	0.92	2.0	ND	1	03/31/06	03/31/06	
Dibromochloromethane	EPA 8260B	6C31009	0.28	1.0	ND	1	03/31/06	03/31/06	
1,2-Dibromoethane (EDB)	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
1,4-Dichlorobenzene	EPA 8260B	6C31009	0.37	1.0	ND	1	03/31/06	03/31/06	
1,2-Dichlorobenzene	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
1,3-Dichlorobenzene	EPA 8260B	6C31009	0.35	1.0	ND	1	03/31/06	03/31/06	
Dichlorodifluoromethane	EPA 8260B	6C31009	0.79	1.0	ND	1	03/31/06	03/31/06	
1,2-Dichloroethane	EPA 8260B	6C31009	0.28	0.50	ND	1	03/31/06	03/31/06	

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 Michele Chamberlin  
 Project Manager

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TAIT Environmental/Boeing  
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 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2080

Sampled: 03/20/06  
 Received: 03/20/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2080-08RE2 (IRZMW002B_WG032006_0001 - Water) - cont.</b>									
Reporting Units: ug/l									
1,1-Dichloroethane	EPA 8260B	6C31009	0.27	1.0	ND	1	03/31/06	03/31/06	
<b>1,1-Dichloroethene</b>	EPA 8260B	6C31009	0.42	1.0	<b>6.2</b>	1	03/31/06	03/31/06	
<b>cis-1,2-Dichloroethene</b>	EPA 8260B	6C31009	0.32	1.0	<b>110</b>	1	03/31/06	03/31/06	
trans-1,2-Dichloroethene	EPA 8260B	6C31009	0.27	1.0	ND	1	03/31/06	03/31/06	
1,2-Dichloropropane	EPA 8260B	6C31009	0.35	1.0	ND	1	03/31/06	03/31/06	
2,2-Dichloropropane	EPA 8260B	6C31009	0.34	1.0	ND	1	03/31/06	03/31/06	
cis-1,3-Dichloropropene	EPA 8260B	6C31009	0.22	0.50	ND	1	03/31/06	03/31/06	
1,1-Dichloropropene	EPA 8260B	6C31009	0.28	1.0	ND	1	03/31/06	03/31/06	
trans-1,3-Dichloropropene	EPA 8260B	6C31009	0.32	0.50	ND	1	03/31/06	03/31/06	
Ethylbenzene	EPA 8260B	6C31009	0.25	1.0	ND	1	03/31/06	03/31/06	
Hexachlorobutadiene	EPA 8260B	6C31009	0.38	1.0	ND	1	03/31/06	03/31/06	
2-Hexanone	EPA 8260B	6C31009	2.6	6.0	ND	1	03/31/06	03/31/06	
Isopropylbenzene	EPA 8260B	6C31009	0.25	1.0	ND	1	03/31/06	03/31/06	
p-Isopropyltoluene	EPA 8260B	6C31009	0.28	1.0	ND	1	03/31/06	03/31/06	
Methyl-tert-butyl Ether (MTBE)	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
Methylene chloride	EPA 8260B	6C31009	0.70	1.0	ND	1	03/31/06	03/31/06	
4-Methyl-2-pentanone (MIBK)	EPA 8260B	6C31009	3.5	5.0	ND	1	03/31/06	03/31/06	
n-Propylbenzene	EPA 8260B	6C31009	0.27	1.0	ND	1	03/31/06	03/31/06	
Styrene	EPA 8260B	6C31009	0.16	1.0	ND	1	03/31/06	03/31/06	
1,1,1,2-Tetrachloroethane	EPA 8260B	6C31009	0.27	1.0	ND	1	03/31/06	03/31/06	
1,1,2,2-Tetrachloroethane	EPA 8260B	6C31009	0.24	1.0	ND	1	03/31/06	03/31/06	
Tetrachloroethene	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
<b>Toluene</b>	EPA 8260B	6C31009	0.36	1.0	<b>1.4</b>	1	03/31/06	03/31/06	
1,2,3-Trichlorobenzene	EPA 8260B	6C31009	0.45	1.0	ND	1	03/31/06	03/31/06	
1,2,4-Trichlorobenzene	EPA 8260B	6C31009	0.48	1.0	ND	1	03/31/06	03/31/06	
1,1,2-Trichloroethane	EPA 8260B	6C31009	0.30	1.0	ND	1	03/31/06	03/31/06	
1,1,1-Trichloroethane	EPA 8260B	6C31009	0.30	1.0	ND	1	03/31/06	03/31/06	
Trichlorofluoromethane	EPA 8260B	6C31009	0.34	2.0	ND	1	03/31/06	03/31/06	
1,2,3-Trichloropropane	EPA 8260B	6C31009	0.40	1.0	ND	1	03/31/06	03/31/06	
1,2,4-Trimethylbenzene	EPA 8260B	6C31009	0.23	1.0	ND	1	03/31/06	03/31/06	
1,3,5-Trimethylbenzene	EPA 8260B	6C31009	0.26	1.0	ND	1	03/31/06	03/31/06	
Vinyl acetate	EPA 8260B	6C31009	1.7	6.0	ND	1	03/31/06	03/31/06	
Vinyl chloride	EPA 8260B	6C31009	0.26	0.50	ND	1	03/31/06	03/31/06	
Xylenes, Total	EPA 8260B	6C31009	0.90	1.0	ND	1	03/31/06	03/31/06	
<i>Surrogate: 4-Bromofluorobenzene (80-120%)</i>							114 %		
<i>Surrogate: Dibromofluoromethane (80-120%)</i>							116 %		
<i>Surrogate: Toluene-d8 (80-120%)</i>							108 %		

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 Michele Chamberlin  
 Project Manager

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Sampled: 03/20/06  
 Received: 03/20/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
<b>Sample ID: IPC2080-09 (IRZMW002B_WG032006_0002 - Water)</b>									
<b>Reporting Units: ug/l</b>									
Iodomethane	EPA 8260B	6C29002	1.0	2.0	ND	1	03/29/06	03/29/06	
Tetrahydrofuran (THF)	EPA 8260B	6C29002	2.3	10	ND	1	03/29/06	03/29/06	
<i>Surrogate: 4-Bromofluorobenzene (80-120%)</i>					96 %				
<i>Surrogate: Dibromofluoromethane (80-120%)</i>					119 %				
<i>Surrogate: Toluene-d8 (80-120%)</i>					101 %				
<b>Sample ID: IPC2080-09RE1 (IRZMW002B_WG032006_0002 - Water)</b>									
<b>Reporting Units: ug/l</b>									
Trichloroethene	EPA 8260B	6C31009	1.3	5.0	330	5	03/31/06	03/31/06	
<i>Surrogate: 4-Bromofluorobenzene (80-120%)</i>					105 %				
<i>Surrogate: Dibromofluoromethane (80-120%)</i>					108 %				
<i>Surrogate: Toluene-d8 (80-120%)</i>					104 %				
<b>Sample ID: IPC2080-09RE2 (IRZMW002B_WG032006_0002 - Water)</b>									
<b>Reporting Units: ug/l</b>									
Acetone	EPA 8260B	6C31009	4.5	10	ND 45	1	03/31/06	03/31/06	
Benzene	EPA 8260B	6C31009	0.28	1.0	ND	1	03/31/06	03/31/06	
Bromobenzene	EPA 8260B	6C31009	0.27	1.0	ND	1	03/31/06	03/31/06	
Bromochloromethane	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
Bromodichloromethane	EPA 8260B	6C31009	0.30	1.0	ND	1	03/31/06	03/31/06	
Bromoform	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
Bromomethane	EPA 8260B	6C31009	0.42	1.0	ND	1	03/31/06	03/31/06	
2-Butanone (MEK)	EPA 8260B	6C31009	3.8	5.0	ND	1	03/31/06	03/31/06	
n-Butylbenzene	EPA 8260B	6C31009	0.37	1.0	ND	1	03/31/06	03/31/06	
sec-Butylbenzene	EPA 8260B	6C31009	0.25	1.0	ND	1	03/31/06	03/31/06	
tert-Butylbenzene	EPA 8260B	6C31009	0.22	1.0	ND	1	03/31/06	03/31/06	
Carbon Disulfide	EPA 8260B	6C31009	0.48	1.0	ND	1	03/31/06	03/31/06	
Carbon tetrachloride	EPA 8260B	6C31009	0.28	0.50	ND	1	03/31/06	03/31/06	
<b>Chlorobenzene</b>	EPA 8260B	6C31009	0.36	1.0	<b>0.48</b>	1	03/31/06	03/31/06	J
Chloroethane	EPA 8260B	6C31009	0.40	2.0	ND	1	03/31/06	03/31/06	
<b>Chloroform</b>	EPA 8260B	6C31009	0.33	1.0	<b>1.3</b>	1	03/31/06	03/31/06	
Chloromethane	EPA 8260B	6C31009	0.30	2.0	ND	1	03/31/06	03/31/06	
2-Chlorotoluene	EPA 8260B	6C31009	0.28	1.0	ND	1	03/31/06	03/31/06	
4-Chlorotoluene	EPA 8260B	6C31009	0.29	1.0	ND	1	03/31/06	03/31/06	
1,2-Dibromo-3-chloropropane	EPA 8260B	6C31009	0.92	2.0	ND	1	03/31/06	03/31/06	
Dibromochloromethane	EPA 8260B	6C31009	0.28	1.0	ND	1	03/31/06	03/31/06	
1,2-Dibromoethane (EDB)	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
1,4-Dichlorobenzene	EPA 8260B	6C31009	0.37	1.0	ND	1	03/31/06	03/31/06	
1,2-Dichlorobenzene	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
1,3-Dichlorobenzene	EPA 8260B	6C31009	0.35	1.0	ND	1	03/31/06	03/31/06	
Dichlorodifluoromethane	EPA 8260B	6C31009	0.79	1.0	ND	1	03/31/06	03/31/06	
1,2-Dichloroethane	EPA 8260B	6C31009	0.28	0.50	ND	1	03/31/06	03/31/06	

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 Michele Chamberlin  
 Project Manager

XQ7ND6

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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2080

Sampled: 03/20/06  
 Received: 03/20/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2080-09RE2 (IRZMW002B_WG032006_0002 - Water) - cont.</b>									
Reporting Units: ug/l									
1,1-Dichloroethane	EPA 8260B	6C31009	0.27	1.0	ND	1	03/31/06	03/31/06	
<b>1,1-Dichloroethene</b>	EPA 8260B	6C31009	0.42	1.0	<b>3.9</b>	1	03/31/06	03/31/06	
<b>cis-1,2-Dichloroethene</b>	EPA 8260B	6C31009	0.32	1.0	<b>94</b>	1	03/31/06	03/31/06	
<b>trans-1,2-Dichloroethene</b>	EPA 8260B	6C31009	0.27	1.0	<b>4.2</b>	1	03/31/06	03/31/06	
1,2-Dichloropropane	EPA 8260B	6C31009	0.35	1.0	ND	1	03/31/06	03/31/06	
2,2-Dichloropropane	EPA 8260B	6C31009	0.34	1.0	ND	1	03/31/06	03/31/06	
cis-1,3-Dichloropropene	EPA 8260B	6C31009	0.22	0.50	ND	1	03/31/06	03/31/06	
1,1-Dichloropropene	EPA 8260B	6C31009	0.28	1.0	ND	1	03/31/06	03/31/06	
trans-1,3-Dichloropropene	EPA 8260B	6C31009	0.32	0.50	ND	1	03/31/06	03/31/06	
Ethylbenzene	EPA 8260B	6C31009	0.25	1.0	ND	1	03/31/06	03/31/06	
Hexachlorobutadiene	EPA 8260B	6C31009	0.38	1.0	ND	1	03/31/06	03/31/06	
2-Hexanone	EPA 8260B	6C31009	2.6	6.0	ND	1	03/31/06	03/31/06	
Isopropylbenzene	EPA 8260B	6C31009	0.25	1.0	ND	1	03/31/06	03/31/06	
p-Isopropyltoluene	EPA 8260B	6C31009	0.28	1.0	ND	1	03/31/06	03/31/06	
Methyl-tert-butyl Ether (MTBE)	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
Methylene chloride	EPA 8260B	6C31009	0.70	1.0	ND	1	03/31/06	03/31/06	
4-Methyl-2-pentanone (MIBK)	EPA 8260B	6C31009	3.5	5.0	ND	1	03/31/06	03/31/06	
n-Propylbenzene	EPA 8260B	6C31009	0.27	1.0	ND	1	03/31/06	03/31/06	
Styrene	EPA 8260B	6C31009	0.16	1.0	ND	1	03/31/06	03/31/06	
1,1,1,2-Tetrachloroethane	EPA 8260B	6C31009	0.27	1.0	ND	1	03/31/06	03/31/06	
1,1,2,2-Tetrachloroethane	EPA 8260B	6C31009	0.24	1.0	ND	1	03/31/06	03/31/06	
Tetrachloroethene	EPA 8260B	6C31009	0.32	1.0	ND	1	03/31/06	03/31/06	
<b>Toluene</b>	EPA 8260B	6C31009	0.36	1.0	<b>1.3</b>	1	03/31/06	03/31/06	
1,2,3-Trichlorobenzene	EPA 8260B	6C31009	0.45	1.0	ND	1	03/31/06	03/31/06	
1,2,4-Trichlorobenzene	EPA 8260B	6C31009	0.48	1.0	ND	1	03/31/06	03/31/06	
1,1,2-Trichloroethane	EPA 8260B	6C31009	0.30	1.0	ND	1	03/31/06	03/31/06	
1,1,1-Trichloroethane	EPA 8260B	6C31009	0.30	1.0	ND	1	03/31/06	03/31/06	
Trichlorofluoromethane	EPA 8260B	6C31009	0.34	2.0	ND	1	03/31/06	03/31/06	
1,2,3-Trichloropropane	EPA 8260B	6C31009	0.40	1.0	ND	1	03/31/06	03/31/06	
1,2,4-Trimethylbenzene	EPA 8260B	6C31009	0.23	1.0	ND	1	03/31/06	03/31/06	
1,3,5-Trimethylbenzene	EPA 8260B	6C31009	0.26	1.0	ND	1	03/31/06	03/31/06	
Vinyl acetate	EPA 8260B	6C31009	1.7	6.0	ND	1	03/31/06	03/31/06	
Vinyl chloride	EPA 8260B	6C31009	0.26	0.50	ND	1	03/31/06	03/31/06	
Xylenes, Total	EPA 8260B	6C31009	0.90	1.0	ND	1	03/31/06	03/31/06	
Surrogate: 4-Bromofluorobenzene (80-120%)					113 %				
Surrogate: Dibromofluoromethane (80-120%)					106 %				
Surrogate: Toluene-d8 (80-120%)					106 %				

**Del Mar Analytical - Irvine**  
 Michele Chamberlin  
 Project Manager

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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2080

Sampled: 03/20/06  
 Received: 03/20/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
<b>Sample ID: IPC2080-10 (IRZB0081_WG032006_0001 - Water)</b>									
<b>Reporting Units: ug/l</b>									
Iodomethane	EPA 8260B	6C29002	1.0	2.0	ND	1	03/29/06	03/29/06	
Tetrahydrofuran (THF)	EPA 8260B	6C29002	2.3	10	ND	1	03/29/06	03/29/06	
<i>Surrogate: 4-Bromofluorobenzene (80-120%)</i>					99 %				
<i>Surrogate: Dibromofluoromethane (80-120%)</i>					121 %				Z2
<i>Surrogate: Toluene-d8 (80-120%)</i>					102 %				
<b>Sample ID: IPC2080-10RE1 (IRZB0081_WG032006_0001 - Water)</b>									
<b>Reporting Units: ug/l</b>									
Vinyl chloride	EPA 8260B	6C31009	5.2	10	1600	20	03/31/06	03/31/06	
<i>Surrogate: 4-Bromofluorobenzene (80-120%)</i>					104 %				
<i>Surrogate: Dibromofluoromethane (80-120%)</i>					112 %				
<i>Surrogate: Toluene-d8 (80-120%)</i>					104 %				
<b>Sample ID: IPC2080-10RE2 (IRZB0081_WG032006_0001 - Water)</b>									
<b>Reporting Units: ug/l</b>									
Acetone	EPA 8260B	6C31009	22	50	ND	5	03/31/06	03/31/06	
Benzene	EPA 8260B	6C31009	1.4	5.0	ND	5	03/31/06	03/31/06	
Bromobenzene	EPA 8260B	6C31009	1.4	5.0	ND	5	03/31/06	03/31/06	
Bromoform	EPA 8260B	6C31009	1.6	5.0	ND	5	03/31/06	03/31/06	
Bromochloromethane	EPA 8260B	6C31009	1.6	5.0	ND	5	03/31/06	03/31/06	
Bromodichloromethane	EPA 8260B	6C31009	1.5	5.0	ND	5	03/31/06	03/31/06	
Chlorobenzene	EPA 8260B	6C31009	1.8	5.0	ND	5	03/31/06	03/31/06	
Chloroform	EPA 8260B	6C31009	1.6	5.0	ND	5	03/31/06	03/31/06	
Chloromethane	EPA 8260B	6C31009	1.5	10	ND	5	03/31/06	03/31/06	
2-Chlorotoluene	EPA 8260B	6C31009	1.4	5.0	ND	5	03/31/06	03/31/06	
2-Butanone (MEK)	EPA 8260B	6C31009	19	25	ND	5	03/31/06	03/31/06	
n-Butylbenzene	EPA 8260B	6C31009	1.8	5.0	ND	5	03/31/06	03/31/06	
sec-Butylbenzene	EPA 8260B	6C31009	1.2	5.0	ND	5	03/31/06	03/31/06	
tert-Butylbenzene	EPA 8260B	6C31009	1.1	5.0	ND	5	03/31/06	03/31/06	
Carbon Disulfide	EPA 8260B	6C31009	2.4	5.0	ND	5	03/31/06	03/31/06	
Carbon tetrachloride	EPA 8260B	6C31009	1.4	2.5	ND	5	03/31/06	03/31/06	
Chloroform	EPA 8260B	6C31009	1.8	5.0	ND	5	03/31/06	03/31/06	
Chloroethane	EPA 8260B	6C31009	2.0	10	ND	5	03/31/06	03/31/06	
Chloromethane	EPA 8260B	6C31009	1.6	5.0	ND	5	03/31/06	03/31/06	
Chlorotoluene	EPA 8260B	6C31009	1.5	10	ND	5	03/31/06	03/31/06	
2-Chlorotoluene	EPA 8260B	6C31009	1.4	5.0	ND	5	03/31/06	03/31/06	
4-Chlorotoluene	EPA 8260B	6C31009	1.4	5.0	ND	5	03/31/06	03/31/06	
1,2-Dibromo-3-chloropropane	EPA 8260B	6C31009	4.6	10	ND	5	03/31/06	03/31/06	
Dibromochloromethane	EPA 8260B	6C31009	1.4	5.0	ND	5	03/31/06	03/31/06	
1,2-Dibromoethane (EDB)	EPA 8260B	6C31009	1.6	5.0	ND	5	03/31/06	03/31/06	
1,4-Dichlorobenzene	EPA 8260B	6C31009	1.8	5.0	ND	5	03/31/06	03/31/06	
1,2-Dichlorobenzene	EPA 8260B	6C31009	1.6	5.0	ND	5	03/31/06	03/31/06	
1,3-Dichlorobenzene	EPA 8260B	6C31009	1.8	5.0	ND	5	03/31/06	03/31/06	
Dichlorodifluoromethane	EPA 8260B	6C31009	4.0	5.0	ND	5	03/31/06	03/31/06	
1,2-Dichloroethane	EPA 8260B	6C31009	1.4	2.5	ND	5	03/31/06	03/31/06	

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TAIT Environmental/Boeing  
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 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2080

Sampled: 03/20/06  
 Received: 03/20/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
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Sample ID: IPC2080-10RE2 (IRZB0081\_WG032006\_0001 - Water) - cont.

Reporting Units: ug/l

1,1-Dichloroethane	EPA 8260B	6C31009	1.4	5.0	ND	5	03/31/06	03/31/06	J
<b>1,1-Dichloroethene</b>	EPA 8260B	6C31009	2.1	5.0	<b>2.2</b>	5	03/31/06	03/31/06	
<b>cis-1,2-Dichloroethene</b>	EPA 8260B	6C31009	1.6	5.0	<b>260</b>	5	03/31/06	03/31/06	
<b>trans-1,2-Dichloroethene</b>	EPA 8260B	6C31009	1.4	5.0	<b>8.1</b>	5	03/31/06	03/31/06	
1,2-Dichloropropane	EPA 8260B	6C31009	1.8	5.0	ND	5	03/31/06	03/31/06	
2,2-Dichloropropane	EPA 8260B	6C31009	1.7	5.0	ND	5	03/31/06	03/31/06	
cis-1,3-Dichloropropene	EPA 8260B	6C31009	1.1	2.5	ND	5	03/31/06	03/31/06	
1,1-Dichloropropene	EPA 8260B	6C31009	1.4	5.0	ND	5	03/31/06	03/31/06	
trans-1,3-Dichloropropene	EPA 8260B	6C31009	1.6	2.5	ND	5	03/31/06	03/31/06	
Ethylbenzene	EPA 8260B	6C31009	1.2	5.0	ND	5	03/31/06	03/31/06	
Hexachlorobutadiene	EPA 8260B	6C31009	1.9	5.0	ND	5	03/31/06	03/31/06	
2-Hexanone	EPA 8260B	6C31009	13	30	ND	5	03/31/06	03/31/06	
Isopropylbenzene	EPA 8260B	6C31009	1.2	5.0	ND	5	03/31/06	03/31/06	
p-Isopropyltoluene	EPA 8260B	6C31009	1.4	5.0	ND	5	03/31/06	03/31/06	
Methyl-tert-butyl Ether (MTBE)	EPA 8260B	6C31009	1.6	5.0	ND	5	03/31/06	03/31/06	
Methylene chloride	EPA 8260B	6C31009	3.5	5.0	ND	5	03/31/06	03/31/06	
4-Methyl-2-pentanone (MIBK)	EPA 8260B	6C31009	18	25	ND	5	03/31/06	03/31/06	
n-Propylbenzene	EPA 8260B	6C31009	1.4	5.0	ND	5	03/31/06	03/31/06	
Styrene	EPA 8260B	6C31009	0.80	5.0	ND	5	03/31/06	03/31/06	
1,1,1,2-Tetrachloroethane	EPA 8260B	6C31009	1.4	5.0	ND	5	03/31/06	03/31/06	
1,1,2,2-Tetrachloroethane	EPA 8260B	6C31009	1.2	5.0	ND	5	03/31/06	03/31/06	
Tetrachloroethene	EPA 8260B	6C31009	1.6	5.0	ND	5	03/31/06	03/31/06	
Toluene	EPA 8260B	6C31009	1.8	5.0	ND	5	03/31/06	03/31/06	
1,2,3-Trichlorobenzene	EPA 8260B	6C31009	2.2	5.0	ND	5	03/31/06	03/31/06	
1,2,4-Trichlorobenzene	EPA 8260B	6C31009	2.4	5.0	ND	5	03/31/06	03/31/06	
1,1,2-Trichloroethane	EPA 8260B	6C31009	1.5	5.0	ND	5	03/31/06	03/31/06	
1,1,1-Trichloroethane	EPA 8260B	6C31009	1.5	5.0	ND	5	03/31/06	03/31/06	
<b>Trichloroethene</b>	EPA 8260B	6C31009	1.3	5.0	<b>36</b>	5	03/31/06	03/31/06	
Trichlorofluoromethane	EPA 8260B	6C31009	1.7	10	ND	5	03/31/06	03/31/06	
1,2,3-Trichloropropane	EPA 8260B	6C31009	2.0	5.0	ND	5	03/31/06	03/31/06	
1,2,4-Trimethylbenzene	EPA 8260B	6C31009	1.2	5.0	ND	5	03/31/06	03/31/06	
1,3,5-Trimethylbenzene	EPA 8260B	6C31009	1.3	5.0	ND	5	03/31/06	03/31/06	
Vinyl acetate	EPA 8260B	6C31009	8.5	30	ND	5	03/31/06	03/31/06	
Xylenes, Total	EPA 8260B	6C31009	4.5	5.0	ND	5	03/31/06	03/31/06	
<i>Surrogate: 4-Bromofluorobenzene (80-120%)</i>					106 %				
<i>Surrogate: Dibromofluoromethane (80-120%)</i>					104 %				
<i>Surrogate: Toluene-d8 (80-120%)</i>					101 %				

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 Michele Chamberlin  
 Project Manager

PC-N75%

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BOE-C6-0050673



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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2080

Sampled: 03/20/06  
 Received: 03/20/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2080-11 (MWB019_WG032006_0001 - Water)</b>									
Reporting Units: ug/l									
Acetone	EPA 8260B	6C30003	4.5	10	ND	1	03/30/06	03/30/06	
Benzene	EPA 8260B	6C30003	0.28	1.0	ND	1	03/30/06	03/30/06	
Bromobenzene	EPA 8260B	6C30003	0.27	1.0	ND	1	03/30/06	03/30/06	
Bromoform	EPA 8260B	6C30003	0.32	1.0	ND	1	03/30/06	03/30/06	
Bromochloromethane	EPA 8260B	6C30003	0.32	1.0	ND	1	03/30/06	03/30/06	
Bromodichloromethane	EPA 8260B	6C30003	0.30	1.0	ND	1	03/30/06	03/30/06	
Bromoform	EPA 8260B	6C30003	0.32	1.0	ND	1	03/30/06	03/30/06	
Bromomethane	EPA 8260B	6C30003	0.42	1.0	ND	1	03/30/06	03/30/06	
2-Butanone (MEK)	EPA 8260B	6C30003	3.8	5.0	ND	1	03/30/06	03/30/06	
n-Butylbenzene	EPA 8260B	6C30003	0.37	1.0	ND	1	03/30/06	03/30/06	M1
sec-Butylbenzene	EPA 8260B	6C30003	0.25	1.0	ND	1	03/30/06	03/30/06	
tert-Butylbenzene	EPA 8260B	6C30003	0.22	1.0	ND	1	03/30/06	03/30/06	
Carbon Disulfide	EPA 8260B	6C30003	0.48	1.0	ND	1	03/30/06	03/30/06	
<b>Carbon tetrachloride</b>	EPA 8260B	6C30003	0.28	0.50	<b>15</b>	1	03/30/06	03/30/06	
Chlorobenzene	EPA 8260B	6C30003	0.36	1.0	ND	1	03/30/06	03/30/06	
Chloroethane	EPA 8260B	6C30003	0.40	2.0	ND	1	03/30/06	03/30/06	
Chloromethane	EPA 8260B	6C30003	0.30	2.0	ND	1	03/30/06	03/30/06	
2-Chlorotoluene	EPA 8260B	6C30003	0.28	1.0	ND	1	03/30/06	03/30/06	
4-Chlorotoluene	EPA 8260B	6C30003	0.29	1.0	ND	1	03/30/06	03/30/06	
1,2-Dibromo-3-chloropropane	EPA 8260B	6C30003	0.92	2.0	ND	1	03/30/06	03/30/06	
Dibromochloromethane	EPA 8260B	6C30003	0.28	1.0	ND	1	03/30/06	03/30/06	
1,2-Dibromoethane (EDB)	EPA 8260B	6C30003	0.32	1.0	ND	1	03/30/06	03/30/06	
1,4-Dichlorobenzene	EPA 8260B	6C30003	0.37	1.0	ND	1	03/30/06	03/30/06	
1,2-Dichlorobenzene	EPA 8260B	6C30003	0.32	1.0	ND	1	03/30/06	03/30/06	
1,3-Dichlorobenzene	EPA 8260B	6C30003	0.35	1.0	ND	1	03/30/06	03/30/06	
Dichlorodifluoromethane	EPA 8260B	6C30003	0.79	1.0	ND	1	03/30/06	03/30/06	
<b>1,2-Dichloroethane</b>	EPA 8260B	6C30003	0.28	0.50	<b>0.40</b>	1	03/30/06	03/30/06	J
<b>1,1-Dichloroethane</b>	EPA 8260B	6C30003	0.27	1.0	<b>0.45</b>	1	03/30/06	03/30/06	J
<b>1,1-Dichloroethene</b>	EPA 8260B	6C30003	0.42	1.0	<b>4.3</b>	1	03/30/06	03/30/06	
cis-1,2-Dichloroethene	EPA 8260B	6C30003	0.32	1.0	ND	1	03/30/06	03/30/06	
trans-1,2-Dichloroethene	EPA 8260B	6C30003	0.27	1.0	ND	1	03/30/06	03/30/06	
1,2-Dichloropropane	EPA 8260B	6C30003	0.35	1.0	ND	1	03/30/06	03/30/06	M1
2,2-Dichloropropane	EPA 8260B	6C30003	0.34	1.0	ND	1	03/30/06	03/30/06	
cis-1,3-Dichloropropene	EPA 8260B	6C30003	0.22	0.50	ND	1	03/30/06	03/30/06	
1,1-Dichloropropene	EPA 8260B	6C30003	0.28	1.0	ND	1	03/30/06	03/30/06	
trans-1,3-Dichloropropene	EPA 8260B	6C30003	0.32	0.50	ND	1	03/30/06	03/30/06	
Ethylbenzene	EPA 8260B	6C30003	0.25	1.0	ND	1	03/30/06	03/30/06	M1
Hexachlorobutadiene	EPA 8260B	6C30003	0.38	1.0	ND	1	03/30/06	03/30/06	M1
2-Hexanone	EPA 8260B	6C30003	2.6	6.0	ND	1	03/30/06	03/30/06	
Iodomethane	EPA 8260B	6C30003	1.0	2.0	ND	1	03/30/06	03/30/06	
Isopropylbenzene	EPA 8260B	6C30003	0.25	1.0	ND	1	03/30/06	03/30/06	
p-Isopropyltoluene	EPA 8260B	6C30003	0.28	1.0	ND	1	03/30/06	03/30/06	

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 Michele Chamberlin  
 Project Manager

*M. Chamberlin*

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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2080

Sampled: 03/20/06  
 Received: 03/20/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
<b>Sample ID: IPC2080-11 (MWB019_WG032006_0001 - Water) - cont.</b>									
Reporting Units: ug/l									
Methyl-tert-butyl Ether (MTBE)	EPA 8260B	6C30003	0.32	1.0	ND	1	03/30/06	03/30/06	
Methylene chloride	EPA 8260B	6C30003	0.70	1.0	ND	1	03/30/06	03/30/06	
4-Methyl-2-pentanone (MIBK)	EPA 8260B	6C30003	3.5	5.0	ND	1	03/30/06	03/30/06	L
n-Propylbenzene	EPA 8260B	6C30003	0.27	1.0	ND	1	03/30/06	03/30/06	
Styrene	EPA 8260B	6C30003	0.16	1.0	ND <i>UJ</i>	1	03/30/06	03/30/06	M2, R-3
1,1,1,2-Tetrachloroethane	EPA 8260B	6C30003	0.27	1.0	ND	1	03/30/06	03/30/06	
1,1,2,2-Tetrachloroethane	EPA 8260B	6C30003	0.24	1.0	ND	1	03/30/06	03/30/06	
Tetrahydrofuran (THF)	EPA 8260B	6C30003	2.3	10	ND	1	03/30/06	03/30/06	
Toluene	EPA 8260B	6C30003	0.36	1.0	ND	1	03/30/06	03/30/06	
1,2,3-Trichlorobenzene	EPA 8260B	6C30003	0.45	1.0	ND	1	03/30/06	03/30/06	
1,2,4-Trichlorobenzene	EPA 8260B	6C30003	0.48	1.0	ND	1	03/30/06	03/30/06	
1,1,2-Trichloroethane	EPA 8260B	6C30003	0.30	1.0	ND	1	03/30/06	03/30/06	
1,1,1-Trichloroethane	EPA 8260B	6C30003	0.30	1.0	ND	1	03/30/06	03/30/06	
Trichlorofluoromethane	EPA 8260B	6C30003	0.34	2.0	ND	1	03/30/06	03/30/06	
1,2,3-Trichloropropane	EPA 8260B	6C30003	0.40	1.0	ND	1	03/30/06	03/30/06	
1,2,4-Trimethylbenzene	EPA 8260B	6C30003	0.23	1.0	ND <i>UJ</i>	1	03/30/06	03/30/06	M2, R-3
1,3,5-Trimethylbenzene	EPA 8260B	6C30003	0.26	1.0	ND	1	03/30/06	03/30/06	
Vinyl acetate	EPA 8260B	6C30003	1.7	6.0	ND	1	03/30/06	03/30/06	
Vinyl chloride	EPA 8260B	6C30003	0.26	0.50	ND	1	03/30/06	03/30/06	
Xylenes, Total	EPA 8260B	6C30003	0.90	1.0	ND	1	03/30/06	03/30/06	
<i>Surrogate: 4-Bromofluorobenzene (80-120%)</i>							96 %		
<i>Surrogate: Dibromofluoromethane (80-120%)</i>							102 %		
<i>Surrogate: Toluene-d8 (80-120%)</i>							106 %		

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 Michele Chamberlin  
 Project Manager

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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2080

Sampled: 03/20/06  
 Received: 03/20/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
<b>Sample ID: IPC2080-11RE1 (MWB019_WG032006_0001 - Water) - cont.</b>									
Reporting Units: ug/l									
Chloroform	EPA 8260B	6C31008	6.6	20	2600	20	03/31/06	03/31/06	
Tetrachloroethene	EPA 8260B	6C31008	6.4	20	190	20	03/31/06	03/31/06	
Trichloroethene	EPA 8260B	6C31008	5.2	20	120	20	03/31/06	03/31/06	
Surrogate: 4-Bromofluorobenzene (80-120%)									
Surrogate: Dibromofluoromethane (80-120%)									
Surrogate: Toluene-d8 (80-120%)									
99 %									
97 %									
98 %									

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 Michele Chamberlin  
 Project Manager

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LDC #: 14879A1  
SDG #: IPC2080  
Laboratory: Del Mar Analytical

## VALIDATION COMPLETENESS WORKSHEET

Tier 1

Date: 5/12/06

Page: /of /

Reviewer: TP  
2nd Reviewer: PD

### METHOD: GC/MS Volatiles (EPA SW 846 Method 8260B)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	△	Sampling dates: 3/20/06
II.	GC/MS Instrument performance check	N	
III.	Initial calibration	N	
IV.	Continuing calibration	N	
V.	Blanks	SW	
VI.	Surrogate spikes	SW	
VII.	Matrix spike/Matrix spike duplicates	SW	
VIII.	Laboratory control samples	SW	LCS
IX.	Regional Quality Assurance and Quality Control	N	
X.	Internal standards	N	
XI.	Target compound identification	N	
XII.	Compound quantitation/CRQLs	N	
XIII.	Tentatively identified compounds (TICs)	N	
XIV.	System performance	N	
XV.	Overall assessment of data	△	
XVI.	Field duplicates	SW	D = 3 + 4      13 + 15      14 + 16
XVII.	Field blanks	N	

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:

*Water*

1	MWB003_WG032006_0001	11	MWB003_WG032006_20001RE	H,S	31	6C29002 ✓
2	MWC011_WG032006_0001	12	MWC011_WG03200622-0001RE		32	6C31009 -
3	IRZMW002B_WG032006_0001	13	IRZMW002B_WG032006-0001RE1		33	6C30003 -
4	IRZMW002B_WG032006_0002	14	IRZMW002B_WG032006-0001RE2		34	6C31008 -
5	IRZB0081_WG032006_0001	15	IRZB0081_WG032006-0002RE1		35	
6	MWB019_WG032006_0001	16	IRZMW002B_WG032006-0002RE2		36	
7	MWB019_WG032006-0001RE1	17	IRZB0081_WG032006-0001RE1		37	
8	MWB019_WG032006.0001RE1	18	IRZB0081_WG032006-0001RE2		38	
9	IRZMW002B_WG032006RE1	19	MWB019_WG032006-0001RE		39	
10	IRZMW002B_WG03200620	20			40	

0002, MSD

RE1

# TARGET COMPOUND WORKSHEET

## METHOD: VOA (EPA SW 846 Method 8260B)

A. Chloromethane*	S. Trichloroethylene	KK. Trichlorofluoromethane	CCC. tert-Butylbenzene	UUU. 1,2-Dichlorotetrafluoroethane
B. Bromomethane	T. Dibromoethylmethane	LL. Methyl- <i>t</i> -butyl ether	DDD. 1,2,4-Trimethylbenzene	VVV. 4-Ethyltoluene
C. Vinyl chloride**	U. 1,1,2-Trichloroethane	MM. 1,2-Dibromo-3-chloropropane	EEE. sec-Butylbenzene	WWW. Ethanol
D. Chloroethane	V. Benzene	NN. Methyl ethyl ketone	FFF. 1,3-Dichlorobenzene	XXX. Di-Isopropyl ether
E. Methylene chloride	W. trans-1,3-Dichloropropene	OO. 2,2-Dichloropropane	GGG. p-Isopropyltoluene	YYY. <i>t</i> -Butanol
F. Acetone	X. Bromoform*	PP. Bromochloromethane	HHH. 1,4-Dichlorobenzene	ZZZ. <i>t</i> -Butyl alcohol
G. Carbon disulfide	Y. 4-Methyl-2-pentanone	QQ. 1,1-Dichloropropene	III. n-Butylbenzene	AAAA. Ethyl <i>t</i> -butyl ether
H. 1,1-Dichloroethene**	Z. 2-Hexanone	RR. Dibromomethane	JJJ. 1,2-Dichlorobenzene	BBBB. <i>t</i> -Amyl methyl ether
I. 1,1-Dichloroethane*	AA. Tetrachloroethene	SS. 1,3-Dichloropropane	KKK. 1,2,4-Trichlorobenzene	CCCC. 1-Chlorohexane
J. 1,2-Dichloroethene, total	BB. 1,1,2,2-Tetrachloroethane*	TT. 1,2-Dibromoethane	LLL. Hexachlorobutadiene	DDDD. Isopropyl alcohol
K. Chloroform**	CC. Toluene**	UU. 1,1,1,2-Tetrachloroethane	MMM. Naphthalene	EEEE. Acetonitrile
L. 1,2-Dichloroethane	DD. Chlorobenzene*	VV. Isopropylbenzene	NNN. 1,2,3-Trichlorobenzene	FFFF. Acrolein
M. 2-Butanone	EE. Ethylbenzene**	WW. Bromobenzene	OOO. 1,3,5-Trichlorobenzene	GGGG. Acrylonitrile
N. 1,1,1-Trichloroethane	FF. Styrene	XX. 1,2,3-Trichloropropane	PPP. trans-1,2-Dichloroethene	HHHH. 1,4-Dioxane
O. Carbon tetrachloride	GG. Xylenes, total	YY. n-Propylbenzene	QQQ. cis-1,2-Dichloroethene	IIII. Isobutyl alcohol
P. Bromodichloromethane	HH. Vinyl acetate	ZZ. 2-Chlorotoluene	RRR. m,p-Xylenes	JJJJ. Methacrylonitrile
Q. 1,2-Dichloropropane**	II. 2-Chloroethylvinyl ether	AAA. 1,3,5-Trimethylbenzene	SSS. o-Xylene	KKKK. Propionitrile
R. cis-1,3-Dichloropropene	JJ. Dichlorodifluoromethane	BBB. 4-Chlorotoluene	TTT. 1,1,2-Trichloro-1,2,2-trifluoroethane	LLLL.

\* = System performance check compounds (SPCC) for RRF ; \*\* = Calibration check compounds (CCC) for %RSD.



LDC #: 14879A  
SDG #: 1PC208C

## **VALIDATION FINDINGS WORKSHEET**

### **Surrogate Spikes**

**METHOD: GC/MS VOA (EPA SW 846 Method 8260B)**

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

<input checked="" type="checkbox"/>	<input type="checkbox"/>	Were all surrogate %R within QC limits?
<input checked="" type="checkbox"/>	<input type="checkbox"/>	If the percent recovery (%R) for one or more surrogates was out of QC limits, was a reanalysis performed to confirm samples with %R out of outside of criteria?

SMC1 (TOL) = Toluene-d8  
 SMC2 (BFB) = Bromofluorobenzene  
 SMC3 (DCE) = 1,2-Dichloroethane  
 SMC4 (DFM) = Dibromofluoromethane

QC Limits (Water)

88-110  
86-115  
80-120  
86-118

QC Limits (Soil)

LDC #: 14879A  
SDG #: PC208C

LDC #: 14879A  
SDG #: PC-2080

## **VALIDATION FINDINGS WORKSHEET**

### **Matrix Spike/Matrix Spike Duplicates**

Page: / of /  
Reviewer: ✓  
2nd Reviewer: ✓

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

Were a matrix spike (MS) and matrix spike duplicate (MSD) analyzed for each matrix in this SDG? If no, indicate which matrix does not have an associated MS/MSD. Soil / Water.

Was a MS/MSD analyzed every 20 samples of each matrix?  
Were the MS/MSD percent recoveries (%R) and the relative  
associated MS/MSD. Soil / Water.

#	Date	MS/MSD ID	Compound	MS %R (Limits)	%R (Limits)	MSD %R (Limits)	RPD (Limits)	Associated Samples	Qualifications
7-8		III	140 (65-135)	( )	( )	( )	( )	C	J/A JUT
		Q	130 (60-125)	( )	( )	( )	( )		
		XXX	145 (60-140)	( )	( )	( )	( )		
		EE	132 (65-130)	( )	( )	( )	( )		
		LL	136 (60-135)	( )	( )	( )	( )		
		S	( )	8 (60-125)	( )	( )	( )		no env spk
		F	( )	23 (45-145)	15 (30)	( )	( )		J/U/J/A
		DD	( )	52 (55-130)	46 (25)	( )	( )		J/U/J/A
		S	( )	48 (40-125)	( )	( )	( )		no env spk
		7-8	( )	( )	( )	( )	( )		
		( )	( )	( )	( )	( )	( )		
		9+10	F	( )	( )	42 (35)	16	J/U/J/A	
		( )	( )	( )	( )	( )	( )		
		( )	( )	( )	( )	( )	( )		
		( )	( )	( )	( )	( )	( )		
		( )	( )	( )	( )	( )	( )		
		( )	( )	( )	( )	( )	( )		
		Compound		QC Limits (Soil)		RPD (Soil)		QC Limits (Water)	RPD (Water)
H.		1,1-Dichloroethene		59-172%		≤ 22%		61-145%	≤ 14%
S.		Trichloroethene		62-137%		≤ 24%		71-120%	≤ 14%
V.		Benzene		66-142%		≤ 21%		76-127%	≤ 11%
CC.		Toluene		59-139%		≤ 21%		76-125%	≤ 13%
DD.		Chlorobenzene		60-133%		≤ 21%		75-130%	≤ 13%

LDC #: 14879A1  
SDG #: 1Pd 2080

# VALIDATION FINDINGS WORKSHEET

## Laboratory Control Samples (LCS)

SDG #: 1PC2080

**METHOD: GC/MS VOA (EPA SW 846 Method 8260B)**

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

Was a LCS required?  
Were the LCS percent  
Y N N/A  
Y N N/A

LDC#: 14879A1  
SDG#: IPC2080

**VALIDATION FINDINGS WORKSHEET**  
**Field Duplicates**

Page: 1 of 1  
Reviewer: PJ  
2nd Reviewer: AC

METHOD: GC/MS VOA(EPA SW846 8260)

- N NA Were field duplicate pairs identified in this SDG?  
 N NA Were target analytes detected in the field duplicate pairs?

Compound	Concentration (ug/L)		RPD	
	13	15		
Trichloroethene	300	330	10	

Compound	Concentration (ug/L)		RPD	
	14	16		
Chlorobenzene	0.42	0.48	13	
Chloroform	1.5	1.3	14	
1,1-Dichloroethene	6.2	3.9	46	
cis-1,2-Dichloroethene	110	94	16	
Toluene	1.4	1.3	7	
trans-1,2-Dichloroethene	1.0u	4.2	123 <i>201</i>	

V:\FIELD DUPLICATES\14879A1.wpd

## **Laboratory Data Consultants, Inc.**

### **Data Validation Report**

**Project/Site Name:** Boeing Realty Corp, Former C-6 Facility, Torrance CA

**Collection Date:** March 21, 2006

**LDC Report Date:** May 15, 2006

**Matrix:** Water

**Parameters:** Volatiles

**Validation Level:** Tier 2

**Laboratory:** Del Mar Analytical

**Sample Delivery Group (SDG):** IPC2190

#### **Sample Identification**

IRZMW001B\_WG032106\_0001  
IRZB0095\_WG032106\_0001  
IRZCMW003B\_WG032106\_0001  
WCC\_6S\_WG032106\_0001  
CMW026\_WG032106\_0001  
IRZMW001B\_WG032106\_0001MS  
IRZMW001B\_WG032106\_0001MSD

## **Introduction**

This data review covers 7 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 8260B for Volatiles.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review (October 1999) as there are no current guidelines for the method stated above.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified a P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blank results are summarized in Section V.

Field duplicates are summarized in Section XVI.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- U     Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
  - J     Indicates an estimated value.
  - R     Quality control indicates the data is not usable.
  - N     Presumptive evidence of presence of the constituent.
  - UJ    Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
  - A     Indicates the finding is based upon technical validation criteria.
  - P     Indicates the finding is related to a protocol/contractual deviation.
- None    Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodices were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. GC/MS Instrument Performance Check**

Instrument performance was checked at 12 hour intervals.

All ion abundance requirements were met.

## **III. Initial Calibration**

Initial calibration was performed using required standard concentrations.

Percent relative standard deviations (%RSD) were less than or equal to 15.0% for each individual compound and less than or equal to 30.0% for calibration check compounds (CCCs).

In the case where %RSD was greater than 15.0%, the laboratory used a calibration curve to evaluate the compound. All coefficients of determination ( $r^2$ ) were greater than or equal to 0.990 .

For the purposes of technical evaluation, all compounds were evaluated against the 30.0% (%RSD) National Functional Guideline criteria. Unless noted above, all compounds were within the validation criteria.

Average relative response factors (RRF) for all volatile target compounds and system performance check compounds (SPCCs) were within method and validation criteria with the following exceptions:

Date	Compound	RRF (Limits)	Associated Samples	Flag	A or P
3/28/06	2-Butanone	0.031 ( $\geq 0.05$ )	IRZMW001B_WG032106_0001 IRZB0095_WG032106_0001 IRZCMW003B_WG032106_0001 WCC_6S_WG032106_0001 CMW026_WG032106_0001 IRZMW001B_WG032106_0001MS IRZMW001B_WG032106_0001MSD 6C29016-BLK	J (all detects) UJ (all non-detects)	A

#### **IV. Continuing Calibration**

Continuing calibration was performed at the required frequencies.

Percent differences (%D) between the initial calibration RRF and the continuing calibration RRF were within the method criteria of less than or equal to 20.0% for calibration check compounds (CCCs).

For the purposes of technical evaluation, all compounds were evaluated against the 25.0% (%D) National Functional Guideline criteria. Unless noted above, all compounds were within the validation criteria.

All of the continuing calibration RRF values were within method and validation criteria with the following exceptions:

Date	Compound	RRF (Limits)	Associated Samples	Flag	A or P
3/29/06	2-Butanone	0.032 ( $\geq 0.05$ )	IRZMW001B_WG032106_0001 IRZB0095_WG032106_0001 IRZCMW003B_WG032106_0001 WCC_6S_WG032106_0001 CMW026_WG032106_0001 IRZMW001B_WG032106_0001MS IRZMW001B_WG032106_0001MSD 6C29016-BLK	J (all detects) UJ (all non-detects)	A

#### **V. Blanks**

Method blanks were reviewed for each matrix as applicable. No volatile contaminants were found in the method blanks with the following exceptions:

Method Blank ID	Analysis Date	Compound TIC (RT in minutes)	Concentration	Associated Samples
6C29016-BLK	3/29/06	Tetrahydrofuran	6.98 ug/L	IRZMW001B_WG032106_0001 IRZB0095_WG032106_0001 IRZCMW003B_WG032106_0001 WCC_6S_WG032106_0001 CMW026_WG032106_0001

Sample concentrations were compared to concentrations detected in the method blanks. The sample concentrations were either not detected or were significantly greater (>10X for common contaminants, >5X for other contaminants) than the concentrations found in the associated method blanks with the following exceptions:

Sample	Compound TIC (RT in minutes)	Reported Concentration	Modified Final Concentration
IRZB0095_WG032106_0001 (10x)	Tetrahydrofuran	64 ug/L	100U ug/L

## **VI. Surrogate Spikes**

Surrogates were added to all samples and blanks as required by the method. All surrogate recoveries (%R) were within QC limits.

## **VII. Matrix Spike/Matrix Spike Duplicates**

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits with the following exceptions:

Spike ID (Associated Samples)	Compound	MS (%R) (Limits)	MSD (%R) (Limits)	RPD (Limits)	Flag	A or P
IRZMW001B_WG032106_0001MS/MSD (IRZMW001B_WG032106_0001)	1,2-Dichloropropane	137 (60-125)	129 (60-125)	-	J (all detects)	A

## **VIII. Laboratory Control Samples (LCS)**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits.

## **IX. Regional Quality Assurance and Quality Control**

Not applicable.

## **X. Internal Standards**

All internal standard areas and retention times were within QC limits.

## **XI. Target Compound Identifications**

Raw data were not reviewed for this SDG.

## **XII. Compound Quantitation and CRQLs**

Raw data were not reviewed for this SDG.

## **XIII. Tentatively Identified Compounds (TICs)**

Raw data were not reviewed for this SDG.

## **XIV. System Performance**

Raw data were not reviewed for this SDG.

## **XV. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **XVI. Field Duplicates**

No field duplicates were identified in this SDG.

## **XVII. Field Blanks**

No field blanks were identified in this SDG.

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Volatiles - Data Qualification Summary - SDG IPC2190**

SDG	Sample	Compound	Flag	A or P	Reason
IPC2190	IRZMW001B_WG032106_0001 IRZB0095_WG032106_0001 IRZCMW003B_WG032106_0001 WCC_6S_WG032106_0001 CMW026_WG032106_0001	2-Butanone	J (all detects) UJ (all non-detects)	A	Initial calibration (RRF)
IPC2190	IRZMW001B_WG032106_0001 IRZB0095_WG032106_0001 IRZCMW003B_WG032106_0001 WCC_6S_WG032106_0001 CMW026_WG032106_0001	2-Butanone	J (all detects) UJ (all non-detects)	A	Continuing calibration (RRF)
IPC2190	IRZMW001B_WG032106_0001	1,2-Dichloropropane	J (all detects)	A	Matrix spike/Matrix spike duplicates (%R)

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Volatiles - Laboratory Blank Data Qualification Summary - SDG IPC2190**

SDG	Sample	Compound TIC (RT in minutes)	Modified Final Concentration	A or P
IPC2190	IRZB0095_WG032106_0001 (10x)	Tetrahydrofuran	100U ug/L	A



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 2520 E. Sunset Rd. #3, Las Vegas, NV 89120 (702) 798-3620 FAX (702) 798-3621

TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2190

Sampled: 03/21/06  
 Received: 03/21/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2190-03 (IRZMW001B_WG032106_0001 - Water)</b>									
Reporting Units: ug/l									
Acetone	EPA 8260B	6C29016	4.5	10	ND	1	03/29/06	03/29/06	
Benzene	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/29/06	
Bromobenzene	EPA 8260B	6C29016	0.27	1.0	ND	1	03/29/06	03/29/06	
Bromoform	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
Bromodichloromethane	EPA 8260B	6C29016	0.30	1.0	ND	1	03/29/06	03/29/06	
Bromoform	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
Bromomethane	EPA 8260B	6C29016	0.42	1.0	ND	1	03/29/06	03/29/06	
2-Butanone (MEK)	EPA 8260B	6C29016	3.8	5.0	ND	1	03/29/06	03/29/06	
n-Butylbenzene	EPA 8260B	6C29016	0.37	1.0	ND	1	03/29/06	03/29/06	
sec-Butylbenzene	EPA 8260B	6C29016	0.25	1.0	ND	1	03/29/06	03/29/06	
tert-Butylbenzene	EPA 8260B	6C29016	0.22	1.0	ND	1	03/29/06	03/29/06	
Carbon Disulfide	EPA 8260B	6C29016	0.48	1.0	ND	1	03/29/06	03/29/06	
Carbon tetrachloride	EPA 8260B	6C29016	0.28	0.50	ND	1	03/29/06	03/29/06	
Chlorobenzene	EPA 8260B	6C29016	0.36	1.0	ND	1	03/29/06	03/29/06	
Chloroethane	EPA 8260B	6C29016	0.40	2.0	ND	1	03/29/06	03/29/06	
<b>Chloroform</b>	EPA 8260B	6C29016	0.33	1.0	<b>2.2</b>	1	03/29/06	03/29/06	
Chloromethane	EPA 8260B	6C29016	0.30	2.0	ND	1	03/29/06	03/29/06	
2-Chlorotoluene	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/29/06	
4-Chlorotoluene	EPA 8260B	6C29016	0.29	1.0	ND	1	03/29/06	03/29/06	
1,2-Dibromo-3-chloropropane	EPA 8260B	6C29016	0.92	2.0	ND	1	03/29/06	03/29/06	
Dibromochloromethane	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/29/06	
1,2-Dibromoethane (EDB)	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
1,4-Dichlorobenzene	EPA 8260B	6C29016	0.37	1.0	ND	1	03/29/06	03/29/06	
1,2-Dichlorobenzene	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
1,3-Dichlorobenzene	EPA 8260B	6C29016	0.35	1.0	ND	1	03/29/06	03/29/06	
Dichlorodifluoromethane	EPA 8260B	6C29016	0.79	1.0	ND	1	03/29/06	03/29/06	
1,2-Dichloroethane	EPA 8260B	6C29016	0.28	0.50	ND	1	03/29/06	03/29/06	
<b>1,1-Dichloroethane</b>	EPA 8260B	6C29016	0.27	1.0	<b>0.65</b>	1	03/29/06	03/29/06	J
<b>1,1-Dichloroethene</b>	EPA 8260B	6C29016	0.42	1.0	<b>31</b>	1	03/29/06	03/29/06	
<b>trans-1,2-Dichloroethene</b>	EPA 8260B	6C29016	0.27	1.0	<b>1.2</b>	1	03/29/06	03/29/06	
1,2-Dichloropropane	EPA 8260B	6C29016	0.35	1.0	ND	1	03/29/06	03/29/06	M1
2,2-Dichloropropane	EPA 8260B	6C29016	0.34	1.0	ND	1	03/29/06	03/29/06	
cis-1,3-Dichloropropene	EPA 8260B	6C29016	0.22	0.50	ND	1	03/29/06	03/29/06	
1,1-Dichloropropene	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/29/06	
trans-1,3-Dichloropropene	EPA 8260B	6C29016	0.32	0.50	ND	1	03/29/06	03/29/06	
Ethylbenzene	EPA 8260B	6C29016	0.25	1.0	ND	1	03/29/06	03/29/06	
Hexachlorobutadiene	EPA 8260B	6C29016	0.38	1.0	ND	1	03/29/06	03/29/06	
2-Hexanone	EPA 8260B	6C29016	2.6	6.0	ND	1	03/29/06	03/29/06	
Iodomethane	EPA 8260B	6C29016	1.0	2.0	ND	1	03/29/06	03/29/06	
Isopropylbenzene	EPA 8260B	6C29016	0.25	1.0	ND	1	03/29/06	03/29/06	
p-Isopropyltoluene	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/29/06	

**Del Mar Analytical - Irvine**

Michele Chamberlin  
 Project Manager

X 25750c

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BOE-C6-0050691



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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2190

Sampled: 03/21/06  
 Received: 03/21/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
<b>Sample ID: IPC2190-03 (IRZMW001B_WG032106_0001 - Water) - cont.</b>									
Reporting Units: ug/l									
Methyl-tert-butyl Ether (MTBE)	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
<b>Methylene chloride</b>	EPA 8260B	6C29016	0.70	1.0	<b>0.77</b>	1	03/29/06	03/29/06	J
4-Methyl-2-pentanone (MIBK)	EPA 8260B	6C29016	3.5	5.0	ND	1	03/29/06	03/29/06	
n-Propylbenzene	EPA 8260B	6C29016	0.27	1.0	ND	1	03/29/06	03/29/06	
Styrene	EPA 8260B	6C29016	0.16	1.0	ND	1	03/29/06	03/29/06	
1,1,1,2-Tetrachloroethane	EPA 8260B	6C29016	0.27	1.0	ND	1	03/29/06	03/29/06	
1,1,2,2-Tetrachloroethane	EPA 8260B	6C29016	0.24	1.0	ND	1	03/29/06	03/29/06	
Tetrachloroethene	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
Tetrahydrofuran (THF)	EPA 8260B	6C29016	2.3	10	ND	1	03/29/06	03/29/06	
Toluene	EPA 8260B	6C29016	0.36	1.0	ND	1	03/29/06	03/29/06	
1,2,3-Trichlorobenzene	EPA 8260B	6C29016	0.45	1.0	ND	1	03/29/06	03/29/06	
1,2,4-Trichlorobenzene	EPA 8260B	6C29016	0.48	1.0	ND	1	03/29/06	03/29/06	
<b>1,1,2-Trichloroethane</b>	EPA 8260B	6C29016	0.30	1.0	<b>0.37</b>	1	03/29/06	03/29/06	J
1,1,1-Trichloroethane	EPA 8260B	6C29016	0.30	1.0	ND	1	03/29/06	03/29/06	
Trichlorofluoromethane	EPA 8260B	6C29016	0.34	2.0	ND	1	03/29/06	03/29/06	
1,2,3-Trichloroproppane	EPA 8260B	6C29016	0.40	1.0	ND	1	03/29/06	03/29/06	
1,2,4-Trimethylbenzene	EPA 8260B	6C29016	0.23	1.0	ND	1	03/29/06	03/29/06	
1,3,5-Trimethylbenzene	EPA 8260B	6C29016	0.26	1.0	ND	1	03/29/06	03/29/06	
Vinyl acetate	EPA 8260B	6C29016	1.7	6.0	ND	1	03/29/06	03/29/06	
Vinyl chloride	EPA 8260B	6C29016	0.26	0.50	ND	1	03/29/06	03/29/06	
Xylenes, Total	EPA 8260B	6C29016	0.90	1.0	ND	1	03/29/06	03/29/06	

Surrogate: 4-Bromofluorobenzene (80-120%)

98 %

Surrogate: Dibromofluoromethane (80-120%)

110 %

Surrogate: Toluene-d8 (80-120%)

107 %

**Del Mar Analytical - Irvine**  
 Michele Chamberlin  
 Project Manager

*LCANR06*

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TAIT Environmental/Boeing  
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 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2190

Sampled: 03/21/06  
 Received: 03/21/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
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Sample ID: IPC2190-03RE1 (IRZMW001B\_WG032106\_0001 - Water) - cont.

Reporting Units: ug/l

cis-1,2-Dichloroethene	EPA 8260B	6C30011	6.4	20	560	20	03/30/06	03/30/06
Trichloroethene	EPA 8260B	6C30011	5.2	20	1200	20	03/30/06	03/30/06
<i>Surrogate: 4-Bromoiodobenzene (80-120%)</i>					104 %			
<i>Surrogate: Dibromoiodomethane (80-120%)</i>					118 %			
<i>Surrogate: Toluene-d8 (80-120%)</i>					106 %			

**Del Mar Analytical - Irvine**  
 Michele Chamberlin  
 Project Manager

X 05/15/06

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Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2190

Sampled: 03/21/06  
 Received: 03/21/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2190-04 (IRZB0095_WG032106_0001 - Water)</b>									
Reporting Units: ug/l									
Acetone	EPA 8260B	6C29016	45	100	ND	10	03/29/06	03/29/06	
Benzene	EPA 8260B	6C29016	2.8	10	ND	10	03/29/06	03/29/06	
Bromobenzene	EPA 8260B	6C29016	2.7	10	ND	10	03/29/06	03/29/06	
Bromoform	EPA 8260B	6C29016	3.2	10	ND	10	03/29/06	03/29/06	
Bromochloromethane	EPA 8260B	6C29016	3.2	10	ND	10	03/29/06	03/29/06	
Bromodichloromethane	EPA 8260B	6C29016	3.0	10	ND	10	03/29/06	03/29/06	
Bromomethane	EPA 8260B	6C29016	3.2	10	ND	10	03/29/06	03/29/06	
2-Butanone (MEK)	EPA 8260B	6C29016	38	50	ND	10	03/29/06	03/29/06	
n-Butylbenzene	EPA 8260B	6C29016	3.7	10	ND	10	03/29/06	03/29/06	
sec-Butylbenzene	EPA 8260B	6C29016	2.5	10	ND	10	03/29/06	03/29/06	
tert-Butylbenzene	EPA 8260B	6C29016	2.2	10	ND	10	03/29/06	03/29/06	
Carbon Disulfide	EPA 8260B	6C29016	4.8	10	ND	10	03/29/06	03/29/06	
Carbon tetrachloride	EPA 8260B	6C29016	2.8	5.0	ND	10	03/29/06	03/29/06	
Chlorobenzene	EPA 8260B	6C29016	3.6	10	ND	10	03/29/06	03/29/06	
Chloroethane	EPA 8260B	6C29016	4.0	20	ND	10	03/29/06	03/29/06	
<b>Chloroform</b>	EPA 8260B	6C29016	3.3	10	<b>7.2</b>	10	03/29/06	03/29/06	J
Chloromethane	EPA 8260B	6C29016	3.0	20	ND	10	03/29/06	03/29/06	
2-Chlorotoluene	EPA 8260B	6C29016	2.8	10	ND	10	03/29/06	03/29/06	
4-Chlorotoluene	EPA 8260B	6C29016	2.9	10	ND	10	03/29/06	03/29/06	
1,2-Dibromo-3-chloropropane	EPA 8260B	6C29016	9.2	20	ND	10	03/29/06	03/29/06	
Dibromochloromethane	EPA 8260B	6C29016	2.8	10	ND	10	03/29/06	03/29/06	
1,2-Dibromoethane (EDB)	EPA 8260B	6C29016	3.2	10	ND	10	03/29/06	03/29/06	
1,4-Dichlorobenzene	EPA 8260B	6C29016	3.7	10	ND	10	03/29/06	03/29/06	
1,2-Dichlorobenzene	EPA 8260B	6C29016	3.2	10	ND	10	03/29/06	03/29/06	
1,3-Dichlorobenzene	EPA 8260B	6C29016	3.5	10	ND	10	03/29/06	03/29/06	
Dichlorodifluoromethane	EPA 8260B	6C29016	7.9	10	ND	10	03/29/06	03/29/06	
1,2-Dichloroethane	EPA 8260B	6C29016	2.8	5.0	ND	10	03/29/06	03/29/06	
1,1-Dichloroethane	EPA 8260B	6C29016	2.7	10	ND	10	03/29/06	03/29/06	
<b>1,1-Dichloroethene</b>	EPA 8260B	6C29016	4.2	10	<b>12</b>	10	03/29/06	03/29/06	
<b>cis-1,2-Dichloroethene</b>	EPA 8260B	6C29016	3.2	10	<b>320</b>	10	03/29/06	03/29/06	
<b>trans-1,2-Dichloroethene</b>	EPA 8260B	6C29016	2.7	10	<b>4.5</b>	10	03/29/06	03/29/06	J
1,2-Dichloropropane	EPA 8260B	6C29016	3.5	10	ND	10	03/29/06	03/29/06	
2,2-Dichloropropane	EPA 8260B	6C29016	3.4	10	ND	10	03/29/06	03/29/06	
cis-1,3-Dichloropropene	EPA 8260B	6C29016	2.2	5.0	ND	10	03/29/06	03/29/06	
1,1-Dichloropropene	EPA 8260B	6C29016	2.8	10	ND	10	03/29/06	03/29/06	
trans-1,3-Dichloropropene	EPA 8260B	6C29016	3.2	5.0	ND	10	03/29/06	03/29/06	
Ethylbenzene	EPA 8260B	6C29016	2.5	10	ND	10	03/29/06	03/29/06	
Hexachlorobutadiene	EPA 8260B	6C29016	3.8	10	ND	10	03/29/06	03/29/06	
2-Hexanone	EPA 8260B	6C29016	26	60	ND	10	03/29/06	03/29/06	
Iodomethane	EPA 8260B	6C29016	10	20	ND	10	03/29/06	03/29/06	
Isopropylbenzene	EPA 8260B	6C29016	2.5	10	ND	10	03/29/06	03/29/06	

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 Michele Chamberlin  
 Project Manager

Karen DC

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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2190

Sampled: 03/21/06  
 Received: 03/21/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
<b>Sample ID: IPC2190-04 (IRZB0095_WG032106_0001 - Water) - cont.</b>									
Reporting Units: ug/l									
p-Isopropyltoluene	EPA 8260B	6C29016	2.8	10	ND	10	03/29/06	03/29/06	
Methyl-tert-butyl Ether (MTBE)	EPA 8260B	6C29016	3.2	10	ND	10	03/29/06	03/29/06	
<b>Methylene chloride</b>	EPA 8260B	6C29016	7.0	10	<b>8.0</b>	10	03/29/06	03/29/06	J
4-Methyl-2-pentanone (MIBK)	EPA 8260B	6C29016	35	50	ND	10	03/29/06	03/29/06	
n-Propylbenzene	EPA 8260B	6C29016	2.7	10	ND	10	03/29/06	03/29/06	
Styrene	EPA 8260B	6C29016	1.6	10	ND	10	03/29/06	03/29/06	
1,1,1,2-Tetrachloroethane	EPA 8260B	6C29016	2.7	10	ND	10	03/29/06	03/29/06	
1,1,2,2-Tetrachloroethane	EPA 8260B	6C29016	2.4	10	ND	10	03/29/06	03/29/06	
<b>Tetrachloroethene</b>	EPA 8260B	6C29016	3.2	10	<b>6.0</b>	10	03/29/06	03/29/06	J
<b>Tetrahydrofuran (THF)</b>	EPA 8260B	6C29016	23	100	<b>64</b> (DDU)	10	03/29/06	03/29/06	B, J
Toluene	EPA 8260B	6C29016	3.6	10	ND	10	03/29/06	03/29/06	
1,2,3-Trichlorobenzene	EPA 8260B	6C29016	4.5	10	ND	10	03/29/06	03/29/06	
1,2,4-Trichlorobenzene	EPA 8260B	6C29016	4.8	10	ND	10	03/29/06	03/29/06	
1,1,2-Trichloroethane	EPA 8260B	6C29016	3.0	10	ND	10	03/29/06	03/29/06	
1,1,1-Trichloroethane	EPA 8260B	6C29016	3.0	10	ND	10	03/29/06	03/29/06	
<b>Trichloroethene</b>	EPA 8260B	6C29016	2.6	10	<b>1100</b>	10	03/29/06	03/29/06	
Trichlorofluoromethane	EPA 8260B	6C29016	3.4	20	ND	10	03/29/06	03/29/06	
1,2,3-Trichloropropane	EPA 8260B	6C29016	4.0	10	ND	10	03/29/06	03/29/06	
1,2,4-Trimethylbenzene	EPA 8260B	6C29016	2.3	10	ND	10	03/29/06	03/29/06	
1,3,5-Trimethylbenzene	EPA 8260B	6C29016	2.6	10	ND	10	03/29/06	03/29/06	
Vinyl acetate	EPA 8260B	6C29016	17	60	ND	10	03/29/06	03/29/06	
<b>Vinyl chloride</b>	EPA 8260B	6C29016	2.6	5.0	<b>590</b>	10	03/29/06	03/29/06	
Xylenes, Total	EPA 8260B	6C29016	9.0	10	ND	10	03/29/06	03/29/06	
<i>Surrogate: 4-Bromofluorobenzene (80-120%)</i>					95 %				
<i>Surrogate: Dibromofluoromethane (80-120%)</i>					108 %				
<i>Surrogate: Toluene-d8 (80-120%)</i>					105 %				

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 Michele Chamberlin  
 Project Manager

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BOE-C6-0050695



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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2190

Sampled: 03/21/06  
 Received: 03/21/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2190-08 (IRZCMW003B_WG032106_0001 - Water)</b>									
Reporting Units: ug/l									
Acetone	EPA 8260B	6C29016	4.5	10	ND	1	03/29/06	03/29/06	
Benzene	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/29/06	
Bromobenzene	EPA 8260B	6C29016	0.27	1.0	ND	1	03/29/06	03/29/06	
Bromoform	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
Bromomethane	EPA 8260B	6C29016	0.30	1.0	ND	1	03/29/06	03/29/06	
2-Butanone (MEK)	EPA 8260B	6C29016	3.8	5.0	ND	1	03/29/06	03/29/06	
n-Butylbenzene	EPA 8260B	6C29016	0.37	1.0	ND	1	03/29/06	03/29/06	
sec-Butylbenzene	EPA 8260B	6C29016	0.25	1.0	ND	1	03/29/06	03/29/06	
tert-Butylbenzene	EPA 8260B	6C29016	0.22	1.0	ND	1	03/29/06	03/29/06	
Carbon Disulfide	EPA 8260B	6C29016	0.48	1.0	ND	1	03/29/06	03/29/06	
Carbon tetrachloride	EPA 8260B	6C29016	0.28	0.50	ND	1	03/29/06	03/29/06	
<b>Chlorobenzene</b>	EPA 8260B	6C29016	0.36	1.0	<b>1.1</b>	1	03/29/06	03/29/06	
Chloroethane	EPA 8260B	6C29016	0.40	2.0	ND	1	03/29/06	03/29/06	
<b>Chloroform</b>	EPA 8260B	6C29016	0.33	1.0	<b>2.1</b>	1	03/29/06	03/29/06	
Chloromethane	EPA 8260B	6C29016	0.30	2.0	ND	1	03/29/06	03/29/06	
2-Chlorotoluene	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/29/06	
4-Chlorotoluene	EPA 8260B	6C29016	0.29	1.0	ND	1	03/29/06	03/29/06	
1,2-Dibromo-3-chloropropane	EPA 8260B	6C29016	0.92	2.0	ND	1	03/29/06	03/29/06	
Dibromochloromethane	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/29/06	
1,2-Dibromoethane (EDB)	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
1,4-Dichlorobenzene	EPA 8260B	6C29016	0.37	1.0	ND	1	03/29/06	03/29/06	
1,2-Dichlorobenzene	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
1,3-Dichlorobenzene	EPA 8260B	6C29016	0.35	1.0	ND	1	03/29/06	03/29/06	
Dichlorodifluoromethane	EPA 8260B	6C29016	0.79	1.0	ND	1	03/29/06	03/29/06	
<b>1,2-Dichloroethane</b>	EPA 8260B	6C29016	0.28	0.50	<b>0.31</b>	1	03/29/06	03/29/06	J
<b>1,1-Dichloroethane</b>	EPA 8260B	6C29016	0.27	1.0	<b>0.86</b>	1	03/29/06	03/29/06	J
<b>1,1-Dichloroethene</b>	EPA 8260B	6C29016	0.42	1.0	<b>.37</b>	1	03/29/06	03/29/06	
<b>trans-1,2-Dichloroethene</b>	EPA 8260B	6C29016	0.27	1.0	<b>2.2</b>	1	03/29/06	03/29/06	
1,2-Dichloropropane	EPA 8260B	6C29016	0.35	1.0	ND	1	03/29/06	03/29/06	
2,2-Dichloropropane	EPA 8260B	6C29016	0.34	1.0	ND	1	03/29/06	03/29/06	
cis-1,3-Dichloropropene	EPA 8260B	6C29016	0.22	0.50	ND	1	03/29/06	03/29/06	
1,1-Dichloropropene	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/29/06	
trans-1,3-Dichloropropene	EPA 8260B	6C29016	0.32	0.50	ND	1	03/29/06	03/29/06	
Ethylbenzene	EPA 8260B	6C29016	0.25	1.0	ND	1	03/29/06	03/29/06	
Hexachlorobutadiene	EPA 8260B	6C29016	0.38	1.0	ND	1	03/29/06	03/29/06	
2-Hexanone	EPA 8260B	6C29016	2.6	6.0	ND	1	03/29/06	03/29/06	
Iodomethane	EPA 8260B	6C29016	1.0	2.0	ND	1	03/29/06	03/29/06	
Isopropylbenzene	EPA 8260B	6C29016	0.25	1.0	ND	1	03/29/06	03/29/06	
p-Isopropyltoluene	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/29/06	

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 Michele Chamberlin  
 Project Manager

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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2190

Sampled: 03/21/06  
 Received: 03/21/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2190-08 (IRZCMW003B_WG032106_0001 - Water) - cont.</b>									
Reporting Units: ug/l									
Methyl-tert-butyl Ether (MTBE)	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
Methylene chloride	EPA 8260B	6C29016	0.70	1.0	ND	1	03/29/06	03/29/06	
4-Methyl-2-pentanone (MIBK)	EPA 8260B	6C29016	3.5	5.0	ND	1	03/29/06	03/29/06	
n-Propylbenzene	EPA 8260B	6C29016	0.27	1.0	ND	1	03/29/06	03/29/06	
Styrene	EPA 8260B	6C29016	0.16	1.0	ND	1	03/29/06	03/29/06	
1,1,1,2-Tetrachloroethane	EPA 8260B	6C29016	0.27	1.0	ND	1	03/29/06	03/29/06	
1,1,2,2-Tetrachloroethane	EPA 8260B	6C29016	0.24	1.0	ND	1	03/29/06	03/29/06	
Tetrachloroethene	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
Tetrahydrofuran (THF)	EPA 8260B	6C29016	2.3	10	ND	1	03/29/06	03/29/06	
Toluene	EPA 8260B	6C29016	0.36	1.0	3.6	1	03/29/06	03/29/06	
1,2,3-Trichlorobenzene	EPA 8260B	6C29016	0.45	1.0	ND	1	03/29/06	03/29/06	
1,2,4-Trichlorobenzene	EPA 8260B	6C29016	0.48	1.0	ND	1	03/29/06	03/29/06	
1,1,2-Trichloroethane	EPA 8260B	6C29016	0.30	1.0	0.41	1	03/29/06	03/29/06	J
1,1,1-Trichloroethane	EPA 8260B	6C29016	0.30	1.0	ND	1	03/29/06	03/29/06	
Trichlorofluoromethane	EPA 8260B	6C29016	0.34	2.0	ND	1	03/29/06	03/29/06	
1,2,3-Trichloropropane	EPA 8260B	6C29016	0.40	1.0	ND	1	03/29/06	03/29/06	
1,2,4-Trimethylbenzene	EPA 8260B	6C29016	0.23	1.0	ND	1	03/29/06	03/29/06	
1,3,5-Trimethylbenzene	EPA 8260B	6C29016	0.26	1.0	ND	1	03/29/06	03/29/06	
Vinyl acetate	EPA 8260B	6C29016	1.7	6.0	ND	1	03/29/06	03/29/06	
Vinyl chloride	EPA 8260B	6C29016	0.26	0.50	3.5	1	03/29/06	03/29/06	
Xylenes, Total	EPA 8260B	6C29016	0.90	1.0	ND	1	03/29/06	03/29/06	

Surrogate: 4-Bromofluorobenzene (80-120%)

99 %

Surrogate: Dibromofluoromethane (80-120%)

114 %

Surrogate: Toluene-d8 (80-120%)

107 %

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 Michele Chamberlin  
 Project Manager

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TAIT Environmental/Boeing  
701 N. Parkcenter Drive  
Santa Ana, CA 92705  
Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
EM2303D/PHASE 01  
Report Number: IPC2190

Sampled: 03/21/06  
Received: 03/21/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2190-08RE1 (IRZCMW003B_WG032106_0001 - Water) - cont.</b>									
Reporting Units: ug/l									
cis-1,2-Dichloroethene	EPA 8260B	6C30026	6.4	20	930	20	03/30/06	03/31/06	
Trichloroethene	EPA 8260B	6C30026	5.2	20	990	20	03/30/06	03/31/06	
Surrogate: 4-Bromofluorobenzene (80-120%)									
Surrogate: Dibromofluoromethane (80-120%)									
Surrogate: Toluene-d8 (80-120%)									
105 %									
119 %									
114 %									

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Michele Chamberlin  
Project Manager

X-057ND

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 2520 E. Sunset Rd. #3, Las Vegas, NV 89120 (702) 798-3620 FAX (702) 798-3621

TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2190

Sampled: 03/21/06  
 Received: 03/21/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2190-09 (WCC_6S_WG032106_0001 - Water)</b>									
Reporting Units: ug/l									
<b>Acetone</b>	EPA 8260B	6C29016	4.5	10	<b>41</b>	1	03/29/06	03/29/06	
<b>Benzene</b>	EPA 8260B	6C29016	0.28	1.0	<b>74</b>	1	03/29/06	03/29/06	
Bromobenzene	EPA 8260B	6C29016	0.27	1.0	ND	1	03/29/06	03/29/06	
Bromoform	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
Bromomethane	EPA 8260B	6C29016	0.30	1.0	ND	1	03/29/06	03/29/06	
Bromochloromethane	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
Bromodichloromethane	EPA 8260B	6C29016	0.30	1.0	ND	1	03/29/06	03/29/06	
Carbon Disulfide	EPA 8260B	6C29016	0.42	1.0	ND	1	03/29/06	03/29/06	
Carbon Tetrachloride	EPA 8260B	6C29016	3.8	5.0	ND <i>45</i>	1	03/29/06	03/29/06	
Chlorobenzene	EPA 8260B	6C29016	0.37	1.0	ND	1	03/29/06	03/29/06	
Chloroethane	EPA 8260B	6C29016	0.25	1.0	ND	1	03/29/06	03/29/06	
Chloroform	EPA 8260B	6C29016	0.22	1.0	ND	1	03/29/06	03/29/06	
Chloromethane	EPA 8260B	6C29016	0.48	1.0	ND	1	03/29/06	03/29/06	
Chlorotoluene	EPA 8260B	6C29016	0.28	0.50	ND	1	03/29/06	03/29/06	
2-Chlorotoluene	EPA 8260B	6C29016	0.36	1.0	ND	1	03/29/06	03/29/06	
4-Chlorotoluene	EPA 8260B	6C29016	0.40	2.0	ND	1	03/29/06	03/29/06	
1,2-Dibromo-3-chloropropane	EPA 8260B	6C29016	0.92	2.0	ND	1	03/29/06	03/29/06	
Dibromochloromethane	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/29/06	
1,2-Dibromoethane (EDB)	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
1,4-Dichlorobenzene	EPA 8260B	6C29016	0.37	1.0	ND	1	03/29/06	03/29/06	
1,2-Dichlorobenzene	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
1,3-Dichlorobenzene	EPA 8260B	6C29016	0.35	1.0	ND	1	03/29/06	03/29/06	
Dichlorodifluoromethane	EPA 8260B	6C29016	0.79	1.0	ND	1	03/29/06	03/29/06	
<b>1,2-Dichloroethane</b>	EPA 8260B	6C29016	0.28	0.50	<b>61</b>	1	03/29/06	03/29/06	
1,2-Dichloropropane	EPA 8260B	6C29016	0.35	1.0	ND	1	03/29/06	03/29/06	
2,2-Dichloropropane	EPA 8260B	6C29016	0.34	1.0	ND	1	03/29/06	03/29/06	
cis-1,3-Dichloropropene	EPA 8260B	6C29016	0.22	0.50	ND	1	03/29/06	03/29/06	
1,1-Dichloropropene	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/29/06	
trans-1,3-Dichloropropene	EPA 8260B	6C29016	0.32	0.50	ND	1	03/29/06	03/29/06	
<b>Ethylbenzene</b>	EPA 8260B	6C29016	0.25	1.0	<b>8.8</b>	1	03/29/06	03/29/06	
Hexachlorobutadiene	EPA 8260B	6C29016	0.38	1.0	ND	1	03/29/06	03/29/06	
2-Hexanone	EPA 8260B	6C29016	2.6	6.0	ND	1	03/29/06	03/29/06	
Iodomethane	EPA 8260B	6C29016	1.0	2.0	ND	1	03/29/06	03/29/06	
<b>Isopropylbenzene</b>	EPA 8260B	6C29016	0.25	1.0	<b>0.39</b>	1	03/29/06	03/29/06	J
p-Isopropyltoluene	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/29/06	
Methyl-tert-butyl Ether (MTBE)	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/29/06	
Methylene chloride	EPA 8260B	6C29016	0.70	1.0	ND	1	03/29/06	03/29/06	
<b>4-Methyl-2-pentanone (MIBK)</b>	EPA 8260B	6C29016	3.5	5.0	<b>4.3</b>	1	03/29/06	03/29/06	J

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 Project Manager

X ACTNDS

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BOE-C6-0050699



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 2520 E. Sunset Rd. #3, Las Vegas, NV 89120 (702) 798-3620 FAX (702) 798-3621

TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2190

Sampled: 03/21/06  
 Received: 03/21/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
<b>Sample ID: IPC2190-09 (WCC_6S_WG032106_0001 - Water) - cont.</b>									
Reporting Units: ug/l									
<b>n-Propylbenzene</b>	EPA 8260B	6C29016	0.27	1.0	<b>0.30</b>	1	03/29/06	03/29/06	J
Styrene	EPA 8260B	6C29016	0.16	1.0	ND	1	03/29/06	03/29/06	
1,1,1,2-Tetrachloroethane	EPA 8260B	6C29016	0.27	1.0	ND	1	03/29/06	03/29/06	
1,1,2,2-Tetrachloroethane	EPA 8260B	6C29016	0.24	1.0	ND	1	03/29/06	03/29/06	
<b>Tetrachloroethene</b>	EPA 8260B	6C29016	0.32	1.0	<b>0.47</b>	1	03/29/06	03/29/06	J
Tetrahydrofuran (THF)	EPA 8260B	6C29016	2.3	10	ND	1	03/29/06	03/29/06	
1,2,3-Trichlorobenzene	EPA 8260B	6C29016	0.45	1.0	ND	1	03/29/06	03/29/06	
1,2,4-Trichlorobenzene	EPA 8260B	6C29016	0.48	1.0	ND	1	03/29/06	03/29/06	
<b>1,1,2-Trichloroethane</b>	EPA 8260B	6C29016	0.30	1.0	<b>46</b>	1	03/29/06	03/29/06	
<b>1,1,1-Trichloroethane</b>	EPA 8260B	6C29016	0.30	1.0	<b>4.0</b>	1	03/29/06	03/29/06	
Trichlorofluoromethane	EPA 8260B	6C29016	0.34	2.0	ND	1	03/29/06	03/29/06	
1,2,3-Trichloropropane	EPA 8260B	6C29016	0.40	1.0	ND	1	03/29/06	03/29/06	
<b>1,2,4-Trimethylbenzene</b>	EPA 8260B	6C29016	0.23	1.0	<b>1.4</b>	1	03/29/06	03/29/06	
<b>1,3,5-Trimethylbenzene</b>	EPA 8260B	6C29016	0.26	1.0	<b>0.61</b>	1	03/29/06	03/29/06	J
Vinyl acetate	EPA 8260B	6C29016	1.7	6.0	ND	1	03/29/06	03/29/06	
<b>Vinyl chloride</b>	EPA 8260B	6C29016	0.26	0.50	<b>5.8</b>	1	03/29/06	03/29/06	
<b>Xylenes, Total</b>	EPA 8260B	6C29016	0.90	1.0	<b>53</b>	1	03/29/06	03/29/06	
Surrogate: 4-Bromo fluorobenzene (80-120%)									
94 %									
Surrogate: Dibromofluoromethane (80-120%)									
109 %									
Surrogate: Toluene-d8 (80-120%)									
108 %									

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 Michele Chamberlin  
 Project Manager

*(Signature)*

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TAIT Environmental/Boeing  
701 N. Parkcenter Drive  
Santa Ana, CA 92705  
Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
EM2303D/PHASE 01  
Report Number: IPC2190

Sampled: 03/21/06  
Received: 03/21/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2190-09RE1 (WCC_6S_WG032106_0001 - Water) - cont.</b>									
Reporting Units: ug/l									
1,1-Dichloroethane	EPA 8260B	6C30011	27	100	310	100	03/30/06	03/30/06	
1,1-Dichloroethene	EPA 8260B	6C30011	42	100	12000	100	03/30/06	03/30/06	
cis-1,2-Dichloroethene	EPA 8260B	6C30011	32	100	5500	100	03/30/06	03/30/06	
trans-1,2-Dichloroethene	EPA 8260B	6C30011	27	100	270	100	03/30/06	03/30/06	
Toluene	EPA 8260B	6C30011	36	100	2600	100	03/30/06	03/30/06	
Trichloroethene	EPA 8260B	6C30011	26	100	1300	100	03/30/06	03/30/06	

Surrogate: 4-Bromofluorobenzene (80-120%)

Surrogate: Dibromofluoromethane (80-120%)

Surrogate: Toluene-d8 (80-120%)

104 %

118 %

106 %

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Michele Chamberlin  
Project Manager

X QTNS06

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 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2190

Sampled: 03/21/06  
 Received: 03/21/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2190-11 (CMW026_WG032106_0001 - Water)</b>									
<b>Reporting Units: ug/l</b>									
Acetone	EPA 8260B	6C29016	4.5	10	ND	1	03/29/06	03/30/06	
Benzene	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/30/06	
Bromobenzene	EPA 8260B	6C29016	0.27	1.0	ND	1	03/29/06	03/30/06	
Bromochloromethane	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/30/06	
Bromodichloromethane	EPA 8260B	6C29016	0.30	1.0	ND	1	03/29/06	03/30/06	
Bromoform	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/30/06	
Bromomethane	EPA 8260B	6C29016	0.42	1.0	ND	1	03/29/06	03/30/06	
2-Butanone (MEK)	EPA 8260B	6C29016	3.8	5.0	ND 4.5	1	03/29/06	03/30/06	
n-Butylbenzene	EPA 8260B	6C29016	0.37	1.0	ND	1	03/29/06	03/30/06	
sec-Butylbenzene	EPA 8260B	6C29016	0.25	1.0	ND	1	03/29/06	03/30/06	
tert-Butylbenzene	EPA 8260B	6C29016	0.22	1.0	ND	1	03/29/06	03/30/06	
Carbon Disulfide	EPA 8260B	6C29016	0.48	1.0	ND	1	03/29/06	03/30/06	
Carbon tetrachloride	EPA 8260B	6C29016	0.28	0.50	ND	1	03/29/06	03/30/06	
Chlorobenzene	EPA 8260B	6C29016	0.36	1.0	ND	1	03/29/06	03/30/06	
Chloroethane	EPA 8260B	6C29016	0.40	2.0	ND	1	03/29/06	03/30/06	
Chloroform	EPA 8260B	6C29016	0.33	1.0	ND	1	03/29/06	03/30/06	
Chloromethane	EPA 8260B	6C29016	0.30	2.0	ND	1	03/29/06	03/30/06	
2-Chlorotoluene	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/30/06	
4-Chlorotoluene	EPA 8260B	6C29016	0.29	1.0	ND	1	03/29/06	03/30/06	
1,2-Dibromo-3-chloropropane	EPA 8260B	6C29016	0.92	2.0	ND	1	03/29/06	03/30/06	
Dibromochloromethane	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/30/06	
1,2-Dibromoethane (EDB)	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/30/06	
1,4-Dichlorobenzene	EPA 8260B	6C29016	0.37	1.0	ND	1	03/29/06	03/30/06	
1,2-Dichlorobenzene	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/30/06	
1,3-Dichlorobenzene	EPA 8260B	6C29016	0.35	1.0	ND	1	03/29/06	03/30/06	
Dichlorodifluoromethane	EPA 8260B	6C29016	0.79	1.0	ND	1	03/29/06	03/30/06	
1,2-Dichloroethane	EPA 8260B	6C29016	0.28	0.50	ND	1	03/29/06	03/30/06	
<b>1,1-Dichloroethane</b>	EPA 8260B	6C29016	0.27	1.0	<b>0.27</b>	1	03/29/06	03/30/06	J
<b>1,1-Dichloroethene</b>	EPA 8260B	6C29016	0.42	1.0	<b>8.4</b>	1	03/29/06	03/30/06	
<b>cis-1,2-Dichloroethene</b>	EPA 8260B	6C29016	0.32	1.0	<b>43</b>	1	03/29/06	03/30/06	
trans-1,2-Dichloroethene	EPA 8260B	6C29016	0.27	1.0	ND	1	03/29/06	03/30/06	
1,2-Dichloropropane	EPA 8260B	6C29016	0.35	1.0	ND	1	03/29/06	03/30/06	
2,2-Dichloropropane	EPA 8260B	6C29016	0.34	1.0	ND	1	03/29/06	03/30/06	
cis-1,3-Dichloropropene	EPA 8260B	6C29016	0.22	0.50	ND	1	03/29/06	03/30/06	
1,1-Dichloropropene	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/30/06	
trans-1,3-Dichloropropene	EPA 8260B	6C29016	0.32	0.50	ND	1	03/29/06	03/30/06	
Ethylbenzene	EPA 8260B	6C29016	0.25	1.0	ND	1	03/29/06	03/30/06	
Hexachlorobutadiene	EPA 8260B	6C29016	0.38	1.0	ND	1	03/29/06	03/30/06	
2-Hexanone	EPA 8260B	6C29016	2.6	6.0	ND	1	03/29/06	03/30/06	
Iodomethane	EPA 8260B	6C29016	1.0	2.0	ND	1	03/29/06	03/30/06	
Isopropylbenzene	EPA 8260B	6C29016	0.25	1.0	ND	1	03/29/06	03/30/06	

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 Michele Chamberlin  
 Project Manager

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 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2190

Sampled: 03/21/06  
 Received: 03/21/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2190-11 (CMW026_WG032106_0001 - Water) - cont.</b>									
Reporting Units: ug/l									
p-Isopropyltoluene	EPA 8260B	6C29016	0.28	1.0	ND	1	03/29/06	03/30/06	
Methyl-tert-butyl Ether (MTBE)	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/30/06	
Methylene chloride	EPA 8260B	6C29016	0.70	1.0	ND	1	03/29/06	03/30/06	
4-Methyl-2-pentanone (MIBK)	EPA 8260B	6C29016	3.5	5.0	ND	1	03/29/06	03/30/06	
n-Propylbenzene	EPA 8260B	6C29016	0.27	1.0	ND	1	03/29/06	03/30/06	
Styrene	EPA 8260B	6C29016	0.16	1.0	ND	1	03/29/06	03/30/06	
1,1,1,2-Tetrachloroethane	EPA 8260B	6C29016	0.27	1.0	ND	1	03/29/06	03/30/06	
1,1,2,2-Tetrachloroethane	EPA 8260B	6C29016	0.24	1.0	ND	1	03/29/06	03/30/06	
Tetrachloroethene	EPA 8260B	6C29016	0.32	1.0	ND	1	03/29/06	03/30/06	
Tetrahydrofuran (THF)	EPA 8260B	6C29016	2.3	10	ND	1	03/29/06	03/30/06	
Toluene	EPA 8260B	6C29016	0.36	1.0	ND	1	03/29/06	03/30/06	
1,2,3-Trichlorobenzene	EPA 8260B	6C29016	0.45	1.0	ND	1	03/29/06	03/30/06	
1,2,4-Trichlorobenzene	EPA 8260B	6C29016	0.48	1.0	ND	1	03/29/06	03/30/06	
1,1,2-Trichloroethane	EPA 8260B	6C29016	0.30	1.0	ND	1	03/29/06	03/30/06	
1,1,1-Trichloroethane	EPA 8260B	6C29016	0.30	1.0	ND	1	03/29/06	03/30/06	
<b>Trichloroethene</b>	EPA 8260B	6C29016	0.26	1.0	<b>12</b>	1	03/29/06	03/30/06	
Trichlorofluoromethane	EPA 8260B	6C29016	0.34	2.0	ND	1	03/29/06	03/30/06	
1,2,3-Trichloropropane	EPA 8260B	6C29016	0.40	1.0	ND	1	03/29/06	03/30/06	
1,2,4-Trimethylbenzene	EPA 8260B	6C29016	0.23	1.0	ND	1	03/29/06	03/30/06	
1,3,5-Trimethylbenzene	EPA 8260B	6C29016	0.26	1.0	ND	1	03/29/06	03/30/06	
Vinyl acetate	EPA 8260B	6C29016	1.7	6.0	ND	1	03/29/06	03/30/06	
Vinyl chloride	EPA 8260B	6C29016	0.26	0.50	ND	1	03/29/06	03/30/06	
Xylenes, Total	EPA 8260B	6C29016	0.90	1.0	ND	1	03/29/06	03/30/06	
<i>Surrogate: 4-Bromofluorobenzene (80-120%)</i>					98 %				
<i>Surrogate: Dibromofluoromethane (80-120%)</i>					112 %				
<i>Surrogate: Toluene-d8 (80-120%)</i>					105 %				

**Del Mar Analytical - Irvine**  
 Michele Chamberlin  
 Project Manager

10/25/2006

LDC #: 14879B1

SDG #: IPC2190

Laboratory: Del Mar Analytical

**VALIDATION COMPLETENESS WORKSHEET**

Tier 2

Date: 5/12/06

Page: 1 of 1

Reviewer: J

2nd Reviewer: d

**METHOD:** GC/MS Volatiles (EPA SW 846 Method 8260B)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	<b>Validation Area</b>		<b>Comments</b>
I.	Technical holding times	A	Sampling dates: 3/21/06
II.	GC/MS Instrument performance check	A	
III.	Initial calibration	SW	% RSD, r <sup>2</sup> 10.990
IV.	Continuing calibration	SW	
V.	Blanks	SW	
VI.	Surrogate spikes	A	
VII.	Matrix spike/Matrix spike duplicates	SW	
VIII.	Laboratory control samples	A	LCS
IX.	Regional Quality Assurance and Quality Control	N	
X.	Internal standards	A	
XI.	Target compound identification	N	
XII.	Compound quantitation/CRQLs	N	
XIII.	Tentatively identified compounds (TICs)	N	
XIV.	System performance	N	
XV.	Overall assessment of data	A	
XVI.	Field duplicates	N	
XVII.	Field blanks	N	

Note: A = Acceptable

ND = No compounds detected

D = Duplicate

N = Not provided/applicable

R = Rinsate

TB = Trip blank

SW = See worksheet

FB = Field blank

EB = Equipment blank

Validated Samples:

Water

✓ 1	IRZMW001B_WG032106_0001	11	6C2901b - BLK	21	3/29	31	
✓ 2	IRZB0095_WG032106_0001	12	6C30011 - BLK	22	3/30	32	
✓ 3	IRZCMW003B_WG032106_0001	13	6C3002b - BLK	23	3/31	33	
✓ 4	WCC_6S_WG032106_0001	14		24		34	
✓ 5	CMW026_WG032106_0001	15		25		35	
6 2	IRZMW001B_WG032106_0001	RE		26		36	
7 3	IRZCMW003B_WG032106_0001	RE		27		37	
8 2	WCC_6S_WG032106_0001	RE		28		38	
9 1	IRZMW001B_WG032106_0001	MS		29		39	
10 1	IRZMW001B_WG032106_0001	MS		30		40	

# TARGET COMPOUND WORKSHEET

## METHOD: VOA (EPA SW 846 Method 8260B)

A. Chloromethane*	S. Trichloroethene	KK. Trichlorofluoromethane	CCC. tert-Butylbenzene	UUU. 1,2-Dichlorotetrafluoroethane
B. Bromomethane	T. Dibromochloromethane	LL. Methyl-1-tert-butyl ether	DDD. 1,2,4-Trimethylbenzene	VVV. 4-Ethyltoluene
C. Vinyl chloride**	U. 1,1,2-Trichloroethane	MM. 1,2-Dibromo-3-chloropropane	EEE. sec-Butylbenzene	WWW. Ethanol
D. Chloroethane	V. Benzene	NN. Methyl ethyl ketone	FFF. 1,3-Dichlorobenzene	XXX. Di-isopropyl ether
E. Methylene chloride	W. trans-1,3-Dichloropropene	OO. 2,2-Dichloropropane	GGG. p-Isopropyltoluene	YYY. tert-Butanol
F. Acetone	X. Bromoform*	PP. Bromochloromethane	HHH. 1,4-Dichlorobenzene	ZZZ. tert-Butyl alcohol
G. Carbon disulfide	Y. 4-Methyl-2-pentanone	QQ. 1,1-Dichloropropane	III. n-Butylbenzene	AAA. Ethyl tert-butyl ether
H. 1,1-Dichloroethene**	Z. 2-Hexanone	RR. Dibromomethane	JJJ. 1,2-Dichlorobenzene	BBB. tert-Amyl methyl ether
I. 1,1-Dichloroethane*	AA. Tetrachloroethane	SS. 1,3-Dichloropropane	KKK. 1,2,4-Trichlorobenzene	CCC. 1-Chlorohexane
J. 1,2-Dichloroethene, total	BB. 1,1,2-Tetrachloroethane*	TT. 1,2-Dibromoethane	LLL. Hexachlorobutadiene	DDD. Isopropyl alcohol
K. Chloroform**	CC. Toluene**	UU. 1,1,1,2-Tetrachloroethane	MMM. Naphthalene	EEE. Acetonitrile
L. 1,2-Dichloroethane	DD. Chlorobenzene*	VV. Isopropylbenzene	NNN. 1,2,3-Trichlorobenzene	FFF. Acrolein
M. 2-Butanone	EE. Ethylbenzene**	WW. Bromobenzene	OOO. 1,3,5-Trichlorobenzene	GGG. Acrylonitrile
N. 1,1,1-Trichloroethane	FF. Styrene	XX. 1,2,3-Trichloropropane	PPP. Trans-1,2-Dichloroethene	HHHH. 1,4-Dioxane
O. Carbon tetrachloride	GG. Xylenes, total	YY. n-Propylbenzene	QQQ. cis-1,2-Dichloroethene	III. Isobutyl alcohol
P. Bromodichloromethane	HH. Vinyl acetate	ZZ. 2-Chlorotoluene	RRR. m,p-Xylenes	JJJJ. Methacrylonitrile
Q. 1,2-Dichloropropane**	II. 2-Chloroethylvinyl ether	AAA. 1,3,5-Trimethylbenzene	SSS. o-Xylene	KKKK. Propionitrile
R. cis-1,3-Dichloropropene	JJ. Dichlorodifluoromethane	BBB. 4-Chlorotoluene	TTT. 1,1,2-Trichloro-1,2,2-trifluoroethane	LLLL.

\* = System performance check compounds (SPCC) for RRF ; \*\* = Calibration check compounds (CCC) for %RSD.

LDC #: 14879B1  
SDG #: PC-2190

# VALIDATION FINDINGS WORKSHEET

## Initial Calibration

**METHOD: GC/MS VOA (EPA SW 846 Method 8260B)**

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

- Did the laboratory perform a 5 point calibration prior to sample analysis?  N  N/A
- Were percent relative standard deviations (%RSD) and relative response factors (RRF) within method criteria for all CCC's and SPCC's?  Y  N/A
- Was a curve fit used for evaluation? If yes, what was the acceptance criteria used for evaluation? \_\_\_\_\_
- Did the initial calibration meet the acceptance criteria?  Y  N/A
- Were all %RSDs and RRFs within the validation criteria of  $\leq 30\% \text{ RSD}$  and  $\geq 0.05 \text{ RRF}$ ?  Y  N/A

Page: 1 of 1  
Reviewer: ✓  
2nd Reviewer: ✓

#	Date	Standard ID	Compound	Finding %RSD (Limit: $\leq 30.0\%$ )	Finding PPF (Limit: $\geq 0.05$ )	Associated Samples	Qualifications
3/28/06	SSC28 VOC	M		0.03	6C29016-BLK	1P 5	J/UJ/A



LDC #: 1487981  
SDG #: 112C2196

# **VALIDATION FINDINGS WORKSHEET**

**Blanks**

Blanks

Page: /  
Reviewer: \_\_\_\_\_  
Reviewer: \_\_\_\_\_

Reviewer: \_\_\_\_\_  
2nd Reviewer: \_\_\_\_\_

**METHOD: GC/MS VOA (EPA SW 846 Method 8260B)**

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

as a method blank associated with every sample in this SDG?

Was a method blank analyzed at least once every 12 hours for each matrix and concentration?

Was there contamination in the method blanks? If yes, what was the source? If no, state every 12 hours for each matrix and concentration?

Blank analysis date: 3/29/06

Conc. units: mg/L

## Associated Samples:

54

Blank analysis date: \_\_\_\_\_

### Conc. units:

All results were qualified using the criteria stated below except those circled

Note: Common contaminants such as Methylene chloride, Acetone, 2-Butanone, Carbon disulfide and TICs that were detected in samples within ten times the associated method blank concentration were qualified as not detected, "U". Other contaminants within five times the method blank concentration were also qualified as not detected, "U".

LDC #: 148793 | SDG #: IPCC21010

## **VALIDATION FINDINGS WORKSHEET**

### **Matrix Spike/Matrix Spike Duplicates**

**METHOD : GC/MS VOA (EPA SW 846 Method 8260B)**

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

<input checked="" type="checkbox"/>	N	N/A	Were a matrix spike (MS) and matrix spike duplicate (MSD) analyzed for each matrix in this SDG? If no, indicate which matrix does not have an associated MS/MSD.
<input checked="" type="checkbox"/>	N	N/A	Was a MS/MSD analyzed every 20 samples of each matrix?
<input checked="" type="checkbox"/>	Y	(N)	Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the QC limits?

## **Laboratory Data Consultants, Inc.**

### **Data Validation Report**

**Project/Site Name:** Boeing Realty Corp, Former C-6 Facility, Torrance CA

**Collection Date:** March 22, 2006

**LDC Report Date:** May 15, 2006

**Matrix:** Water

**Parameters:** Volatiles

**Validation Level:** Tier 3

**Laboratory:** Del Mar Analytical

**Sample Delivery Group (SDG):** IPC2080

#### **Sample Identification**

CMW0002\_WG032206\_0001

## **Introduction**

This data review covers one water sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 8260B for Volatiles.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review (October 1999) as there are no current guidelines for the method stated above.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified a P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blank results are summarized in Section V.

Field duplicates are summarized in Section XVI.

The following are definitions of the data qualifiers:

- U     Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- J     Indicates an estimated value.
- R     Quality control indicates the data is not usable.
- N     Presumptive evidence of presence of the constituent.
- UJ    Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A     Indicates the finding is based upon technical validation criteria.
- P     Indicates the finding is related to a protocol/contractual deviation.
- None    Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. GC/MS Instrument Performance Check**

Instrument performance was checked at 12 hour intervals.

All ion abundance requirements were met.

## **III. Initial Calibration**

Initial calibration was performed using required standard concentrations.

Percent relative standard deviations (%RSD) were less than or equal to 15.0% for each individual compound and less than or equal to 30.0% for calibration check compounds (CCCs).

In the case where %RSD was greater than 15.0%, the laboratory used a calibration curve to evaluate the compound. All coefficients of determination ( $r^2$ ) were greater than or equal to 0.990 .

For the purposes of technical evaluation, all compounds were evaluated against the 30.0% (%RSD) National Functional Guideline criteria. Unless noted above, all compounds were within the validation criteria.

Average relative response factors (RRF) for all volatile target compounds and system performance check compounds (SPCCs) were within method and validation criteria with the following exceptions:

Date	Compound	RRF (Limits)	Associated Samples	Flag	A or P
3/26/06	2-Butanone	0.040 ( $\geq 0.05$ )	CMW0002_WG032206_0001 6D03015-BLK1	J (all detects) UJ (all non-detects)	A

## **IV. Continuing Calibration**

Continuing calibration was performed at the required frequencies.

Percent differences (%D) between the initial calibration RRF and the continuing calibration RRF were within the method criteria of less than or equal to 20.0% for calibration check compounds (CCCs).

For the purposes of technical evaluation, all compounds were evaluated against the 25.0% (%D) National Functional Guideline criteria. Unless noted above, all compounds were within the validation criteria.

All of the continuing calibration RRF values were within method and validation criteria with the following exceptions:

Date	Compound	RRF (Limits)	Associated Samples	Flag	A or P
4/3/06	2-Butanone	0.040 ( $\geq$ 0.05)	CMW0002_WG032206_0001 6D03015-BLK1	J (all detects) UJ (all non-detects)	A

## V. Blanks

Method blanks were reviewed for each matrix as applicable. No volatile contaminants were found in the method blanks with the following exceptions:

Method Blank ID	Analysis Date	Compound TIC (RT in minutes)	Concentration	Associated Samples
6D03015-BLK	4/3/06	Tetrahydrofuran	3.42 ug/L	CMW0002_WG032206_0001

Sample concentrations were compared to concentrations detected in the method blanks. The sample concentrations were either not detected or were significantly greater (>10X for common contaminants, >5X for other contaminants) than the concentrations found in the associated method blanks with the following exceptions:

Sample	Compound TIC (RT in minutes)	Reported Concentration	Modified Final Concentration
CMW0002_WG032206_0001 (2x)	Tetrahydrofuran	6.1 ug/L	20U ug/L

## VI. Surrogate Spikes

Surrogates were added to all samples and blanks as required by the method. All surrogate recoveries (%R) were within QC limits with the following exceptions:

Sample	Surrogate	%R (Limits)	Compound	Flag	A or P
CMW0002_WG032206_0001	Bromofluorobenzene	128 (80-120)	All TCL compounds except Chlorobenzene	J (all detects)	A

## **VII. Matrix Spike/Matrix Spike Duplicates**

The laboratory has indicated that there were no matrix spike (MS) and matrix spike duplicate (MSD) analyses specified for the samples in this SDG, and therefore matrix spike and matrix spike duplicate analyses were not performed for this SDG.

## **VIII. Laboratory Control Samples (LCS)**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits.

## **IX. Regional Quality Assurance and Quality Control**

Not applicable.

## **X. Internal Standards**

All internal standard areas and retention times were within QC limits

## **XI. Target Compound Identifications**

All target compound identifications were within validation criteria.

## **XII. Compound Quantitation and CRQLs**

All compound quantitation and CRQLs were within validation criteria.

## **XIII. Tentatively Identified Compounds (TICs)**

Tentatively identified compounds were not reported by the laboratory.

## **XIV. System Performance**

The system performance was acceptable.

## **XV. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **XVI. Field Duplicates**

No field duplicates were identified in this SDG.

## **XVII. Field Blanks**

No field blanks were identified in this SDG.

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Volatiles - Data Qualification Summary - SDG IPC2080**

SDG	Sample	Compound	Flag	A or P	Reason
IPC2080	CMW0002_WG032206_0001	2-Butanone	J (all detects) UJ (all non-detects)	A	Initial calibration (RRF)
IPC2080	CMW0002_WG032206_0001	2-Butanone	J (all detects) UJ (all non-detects)	A	Continuing calibration (RRF)
IPC2080	CMW0002_WG032206_0001	All TCL compounds except Chlorobenzene	J (all detects)	A	Surrogate recovery (%R)

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Volatiles - Laboratory Blank Data Qualification Summary - SDG IPC2080**

SDG	Sample	Compound TIC (RT in minutes)	Modified Final Concentration	A or P
IPC2080	CMW0002_WG032206_0001 (2x)	Tetrahydrofuran	20U ug/L	A



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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2325

Sampled: 03/22/06  
 Received: 03/22/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2325-09 (CMW0002_WG032206_0001 - Water)</b>									
Reporting Units: ug/l									
Acetone	EPA 8260B	6D03015	9.0	20	ND	2	04/03/06	04/03/06	
<b>Benzene</b>	EPA 8260B	6D03015	0.56	2.0	70 J	2	04/03/06	04/03/06	
Bromobenzene	EPA 8260B	6D03015	0.54	2.0	ND	2	04/03/06	04/03/06	
Bromoform	EPA 8260B	6D03015	0.64	2.0	ND	2	04/03/06	04/03/06	
Bromomethane	EPA 8260B	6D03015	0.84	2.0	ND	2	04/03/06	04/03/06	
2-Butanone (MEK)	EPA 8260B	6D03015	7.6	10	ND UJ	2	04/03/06	04/03/06	
n-Butylbenzene	EPA 8260B	6D03015	0.74	2.0	ND	2	04/03/06	04/03/06	
sec-Butylbenzene	EPA 8260B	6D03015	0.50	2.0	ND	2	04/03/06	04/03/06	
tert-Butylbenzene	EPA 8260B	6D03015	0.44	2.0	ND	2	04/03/06	04/03/06	
Carbon Disulfide	EPA 8260B	6D03015	0.96	2.0	ND	2	04/03/06	04/03/06	
Carbon tetrachloride	EPA 8260B	6D03015	0.56	1.0	ND	2	04/03/06	04/03/06	
Chloroethane	EPA 8260B	6D03015	0.80	4.0	ND	2	04/03/06	04/03/06	
<b>Chloroform</b>	EPA 8260B	6D03015	0.66	2.0	1.2 J	2	04/03/06	04/03/06	J
Chloromethane	EPA 8260B	6D03015	0.60	4.0	ND	2	04/03/06	04/03/06	
2-Chlorotoluene	EPA 8260B	6D03015	0.56	2.0	ND	2	04/03/06	04/03/06	
4-Chlorotoluene	EPA 8260B	6D03015	0.58	2.0	ND	2	04/03/06	04/03/06	
1,2-Dibromo-3-chloropropane	EPA 8260B	6D03015	1.8	4.0	ND	2	04/03/06	04/03/06	
Dibromochloromethane	EPA 8260B	6D03015	0.56	2.0	ND	2	04/03/06	04/03/06	
1,2-Dibromoethane (EDB)	EPA 8260B	6D03015	0.64	2.0	ND	2	04/03/06	04/03/06	
<b>1,4-Dichlorobenzene</b>	EPA 8260B	6D03015	0.74	2.0	8.9 J	2	04/03/06	04/03/06	
<b>1,2-Dichlorobenzene</b>	EPA 8260B	6D03015	0.64	2.0	1.4 J	2	04/03/06	04/03/06	J
1,3-Dichlorobenzene	EPA 8260B	6D03015	0.70	2.0	ND	2	04/03/06	04/03/06	
Dichlorodifluoromethane	EPA 8260B	6D03015	1.6	2.0	ND	2	04/03/06	04/03/06	
1,2-Dichloroethane	EPA 8260B	6D03015	0.56	1.0	ND	2	04/03/06	04/03/06	
1,1-Dichloroethane	EPA 8260B	6D03015	0.54	2.0	ND	2	04/03/06	04/03/06	
<b>1,1-Dichloroethene</b>	EPA 8260B	6D03015	0.84	2.0	1.5 J	2	04/03/06	04/03/06	J
<b>cis-1,2-Dichloroethene</b>	EPA 8260B	6D03015	0.64	2.0	11 J	2	04/03/06	04/03/06	
trans-1,2-Dichloroethene	EPA 8260B	6D03015	0.54	2.0	ND	2	04/03/06	04/03/06	
1,2-Dichloropropane	EPA 8260B	6D03015	0.70	2.0	ND	2	04/03/06	04/03/06	
2,2-Dichloropropane	EPA 8260B	6D03015	0.68	2.0	ND	2	04/03/06	04/03/06	
cis-1,3-Dichloropropene	EPA 8260B	6D03015	0.44	1.0	ND	2	04/03/06	04/03/06	
1,1-Dichloropropene	EPA 8260B	6D03015	0.56	2.0	ND	2	04/03/06	04/03/06	
trans-1,3-Dichloropropene	EPA 8260B	6D03015	0.64	1.0	ND	2	04/03/06	04/03/06	
<b>Ethylbenzene</b>	EPA 8260B	6D03015	0.50	2.0	1.3 J	2	04/03/06	04/03/06	
Hexachlorobutadiene	EPA 8260B	6D03015	0.76	2.0	ND	2	04/03/06	04/03/06	
2-Hexanone	EPA 8260B	6D03015	5.2	12	ND	2	04/03/06	04/03/06	
Iodomethane	EPA 8260B	6D03015	2.0	4.0	ND	2	04/03/06	04/03/06	
Isopropylbenzene	EPA 8260B	6D03015	0.50	2.0	ND	2	04/03/06	04/03/06	
p-Isopropyltoluene	EPA 8260B	6D03015	0.56	2.0	ND	2	04/03/06	04/03/06	

**Del Mar Analytical - Irvine**  
 Michele Chamberlin  
 Project Manager

*Logfile*

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 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2325

Sampled: 03/22/06  
 Received: 03/22/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2325-09 (CMW0002_WG032206_0001 - Water) - cont.</b>									
Reporting Units: ug/l									
Methyl-tert-butyl Ether (MTBE)	EPA 8260B	6D03015	0.64	2.0	ND	2	04/03/06	04/03/06	
Methylene chloride	EPA 8260B	6D03015	1.4	2.0	ND	2	04/03/06	04/03/06	
4-Methyl-2-pentanone (MIBK)	EPA 8260B	6D03015	7.0	10	ND	2	04/03/06	04/03/06	
n-Propylbenzene	EPA 8260B	6D03015	0.54	2.0	ND	2	04/03/06	04/03/06	
Styrene	EPA 8260B	6D03015	0.32	2.0	ND	2	04/03/06	04/03/06	
1,1,1,2-Tetrachloroethane	EPA 8260B	6D03015	0.54	2.0	ND	2	04/03/06	04/03/06	
1,1,2,2-Tetrachloroethane	EPA 8260B	6D03015	0.48	2.0	ND	2	04/03/06	04/03/06	
<b>Tetrachloroethene</b>	EPA 8260B	6D03015	0.64	2.0	3.2 J	2	04/03/06	04/03/06	
<b>Tetrahydrofuran (THF)</b>	EPA 8260B	6D03015	4.6	20	6.1 20 U.S.	2	04/03/06	04/03/06	B, J
Toluene	EPA 8260B	6D03015	0.72	2.0	ND	2	04/03/06	04/03/06	
1,2,3-Trichlorobenzene	EPA 8260B	6D03015	0.90	2.0	ND	2	04/03/06	04/03/06	
1,2,4-Trichlorobenzene	EPA 8260B	6D03015	0.96	2.0	ND	2	04/03/06	04/03/06	
1,1,2-Trichloroethane	EPA 8260B	6D03015	0.60	2.0	ND	2	04/03/06	04/03/06	
1,1,1-Trichloroethane	EPA 8260B	6D03015	0.60	2.0	ND	2	04/03/06	04/03/06	
<b>Trichloroethene</b>	EPA 8260B	6D03015	0.52	2.0	330 J	2	04/03/06	04/03/06	
Trichlorofluoromethane	EPA 8260B	6D03015	0.68	4.0	ND	2	04/03/06	04/03/06	
1,2,3-Trichloropropane	EPA 8260B	6D03015	0.80	2.0	ND	2	04/03/06	04/03/06	
1,2,4-Trimethylbenzene	EPA 8260B	6D03015	0.46	2.0	ND	2	04/03/06	04/03/06	
1,3,5-Trimethylbenzene	EPA 8260B	6D03015	0.52	2.0	ND	2	04/03/06	04/03/06	
Vinyl acetate	EPA 8260B	6D03015	3.4	12	ND	2	04/03/06	04/03/06	
Vinyl chloride	EPA 8260B	6D03015	0.52	1.0	ND	2	04/03/06	04/03/06	
<b>Xylenes, Total</b>	EPA 8260B	6D03015	1.8	2.0	2.3 J	2	04/03/06	04/03/06	
<i>Surrogate: 4-Bromofluorobenzene (80-120%)</i>					128 %				A-01, ZX
<i>Surrogate: Dibromofluoromethane (80-120%)</i>					111 %				
<i>Surrogate: Toluene-d8 (80-120%)</i>					101 %				

**Del Mar Analytical - Irvine**  
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 Project Manager

*MLDTKWD*

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Project ID: Boeing C-6 Torrance  
EM2303D/PHASE 01  
Report Number: IPC2325

Sampled: 03/22/06  
Received: 03/22/06

### VOLATILE ORGANICS by GC/MS (EPA 5030B/8260B)

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2325-09RE1 (CMW0002_WG032206_0001 - Water) - cont.</b>									
Reporting Units: ug/l									
Chlorobenzene	EPA 8260B	6D01007	18	50	6900	50	04/01/06	04/01/06	
Surrogate: 4-Bromofluorobenzene (80-120%)					104 %				
Surrogate: Dibromofluoromethane (80-120%)					112 %				
Surrogate: Toluene-d8 (80-120%)					103 %				

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Project Manager

L047576

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IPC2325 <Page 25 of 73>

LDC #: 14879C1

SDG #: IPC2325

Laboratory: Del Mar Analytical

## VALIDATION COMPLETENESS WORKSHEET

Tier 3

Date: 5/15/06

Page: 1 of 1

Reviewer: P

2nd Reviewer: LC

METHOD: GC/MS Volatiles (EPA SW 846 Method 8260B)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	△	Sampling dates: 3/22/06
II.	GC/MS Instrument performance check	△	
III.	Initial calibration	SW	% RSD, $\sigma^2 = 10.990$
IV.	Continuing calibration	SW	
V.	Blanks	SW	
VI.	Surrogate spikes	SW	
VII.	Matrix spike/Matrix spike duplicates	N	client specified
VIII.	Laboratory control samples	A	LCS
IX.	Regional Quality Assurance and Quality Control	N	
X.	Internal standards	A	
XI.	Target compound identification	△	
XII.	Compound quantitation/CRQLs	△	
XIII.	Tentatively identified compounds (TICs)	N	not reported
XIV.	System performance	A	
XV.	Overall assessment of data	A	
XVI.	Field duplicates	N	
XVII.	Field blanks	N	

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet

ND = No compounds detected  
 R = Rinsate  
 FB = Field blank

D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples:

1	CMW0002_WG032206_0001	11	6D03015 - BLK1	21		31	
2	CMW0002_WG032206_0001	12	6D01007 - BLK1	22	for DD only	32	
3		13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

# TARGET COMPOUND WORKSHEET

## METHOD: VOA (EPA SW 846 Method 8260B)

A. Chloromethane*	S. Trichloroethene	KK. Trichlorofluoromethane	CCC. tert-Butylbenzene	UUU. 1,2-Dichlorotetrafluoroethane
B. Bromomethane	T. Dibromochloromethane	LL. Methyl-tert-Butyl ether	DDD. 1,2,4-Trimethylbenzene	VVV. 4-Ethyltoluene
C. Vinyl chloride**	U. 1,1,2-Trichloroethane	MM. 1,2-Dibromo-3-chloropropane	EEE. sec-Butylbenzene	WWW. Ethanol
D. Chloroethane	V. Benzene	NN. Methyl ethyl ketone	FFF. 1,3-Dichlorobenzene	XXX. Di-Isopropyl ether
E. Methylene chloride	W. trans-1,3-Dichloropropene	OO. 2,2-Dichloropropane	GGG. p-Isopropyltoluene	YYY. tert-Butanol
F. Acetone	X. Bromoform*	PP. Bromochloromethane	HHH. 1,4-Dichlorobenzene	ZZZ. tert-Butyl alcohol
G. Carbon disulfide	Y. 4-Methyl-2-pentanone	QQ. 1,1-Dichloropropane	III. n-Butylbenzene	AAAA. Ethyl tert-butyl ether
H. 1,1-Dichloroethene**	Z. 2-Hexanone	RR. Dibromomethane	JJJ. 1,2-Dichlorobenzene	BBBB. tert-Amyl methyl ether
I. 1,1-Dichloroethane*	AA. Tetrachloroethene	SS. 1,3-Dichloropropane	KKK. 1,2,4-Trichlorobenzene	CCCC. 1-Chlorohexane
J. 1,2-Dichloroethene, total	BB. 1,1,2,2-Tetrachloroethane*	TT. 1,2-Dibromoethane	LLL. Hexachlorobutadiene	DDDD. Isopropyl alcohol
K. Chloroform**	CC. Toluene**	UU. 1,1,1,2-Tetrachloroethane	MM. Naphthalene	EEEE. Acetonitrile
L. 1,2-Dichloroethane	DD. Chlorobenzene*	VV. Isopropylbenzene	NNN. 1,2,3-Trichlorobenzene	FFFF. Acrolein
M. 2-Butanone	EE. Ethylbenzene**	WW. Bromobenzene	OOO. 1,3,5-Trichlorobenzene	GGGG. Acrylonitrile
N. 1,1,1-Trichloroethane	FF. Styrene	XX. 1,2,3-Trichloropropane	PPP. trans-1,2-Dichloroethene	HHHH. 1,4-Dioxane
O. Carbon tetrachloride	GG. Xylenes, total	YY. n-Propylbenzene	QQQ. cis-1,2-Dichloroethene	IIII. Isobutyl alcohol
P. Bromodichloromethane	HH. Vinyl acetate	ZZ. 2-Chlorotoluene	RRR. m,p-Xylenes	JJJJ. Methacrylonitrile
Q. 1,2-Dichloropropane**	II. 2-Chloroethylvinyl ether	AAA. 1,3,5-Trimethylbenzene	SSS. o-Xylene	KKKK. Propionitrile
R. dis-1,3-Dichloropropene	JJ. Dichlorodifluoromethane	BBB. 4-Chlorotoluene	TTT. 1,1,2-Trichloro-1,2,2-trifluoroethane	LLLL.

\* = System performance check compounds (SPCC) for RRF ; \*\* = Calibration check compounds (CCC) for %RSD.

LDC #: 148794  
SDG #: 1PC232

## VALIDATION FINDINGS CHECKLIST

Page: 1 of 2  
Reviewer: BS  
2nd Reviewer: BL

**Method:** Volatiles (EPA SW 846 Method 8260B)

Validation Area	Yes	No	NA	Findings/Comments
<b>I. Technical holding times</b>				
All technical holding times were met.	/			
Cooler temperature criteria was met.	/			
<b>II. GC/MS instrument performance check</b>				
Were the BFB performance results reviewed and found to be within the specified criteria?	/			
Were all samples analyzed within the 12 hour clock criteria?	/			
<b>III. Initial calibration</b>				
Did the laboratory perform a 5 point calibration prior to sample analysis?	/			
Were all percent relative standard deviations (%RSD) and relative response factors (RRF) within method criteria for all CCCs and SPCCs?	/			
Was a curve fit used for evaluation?	/			
Did the initial calibration meet the curve fit acceptance criteria of > 0.990?	/			
Were all percent relative standard deviations (%RSD) $\leq$ 30% and relative response factors (RRF) $>$ 0.05?		/		
<b>IV. Continuing calibration</b>				
Was a continuing calibration standard analyzed at least once every 12 hours for each instrument?	/			
Were all percent differences (%D) and relative response factors (RRF) within method criteria for all CCCs and SPCCs?	/			
Were all percent differences (%D) $\leq$ 25% and relative response factors (RRF) $\geq$ 0.05?		/		
<b>V. Blanks</b>				
Was a method blank associated with every sample in this SDG?	/			
Was a method blank analyzed at least once every 12 hours for each matrix and concentration?	/			
Was there contamination in the method blanks? If yes, please see the Blanks validation completeness worksheet.		/		
<b>VI. Surrogate Spikes</b>				
Were all surrogate %R within QC limits?		/		
If the percent recovery (%R) for one or more surrogates was out of QC limits, was a reanalysis performed to confirm samples with %R outside of criteria?	/			
<b>VII. Matrix spike/Matrix spike duplicates</b>				
Were a matrix spike (MS) and matrix spike duplicate (MSD) analyzed for each matrix in this SDG? If no, indicate which matrix does not have an associated MS/MSD. Soil / Water.			/	
Was a MS/MSD analyzed every 20 samples of each matrix?			/	
Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the QC limits?			/	
<b>VIII. Laboratory control samples</b>				
Was an LCS analyzed for this SDG?	/			

LDC #: 11879C1  
SDG #: 1PC2325

### VALIDATION FINDINGS CHECKLIST

Page: 2 of 2  
Reviewer: B  
2nd Reviewer: A

Validation Area	Yes	No	NA	Findings/Comments
Was an LCS analyzed per analytical batch?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were the LCS percent recoveries (%R) and relative percent difference (RPD) within the QC limits?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>X. Regional Quality Assurance and Quality Control</b>				
Were performance evaluation (PE) samples performed?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
Were the performance evaluation (PE) samples within the acceptance limits?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	
<b>XI. Internal standards</b>				
Were internal standard area counts within -50% or +100% of the associated calibration standard?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were retention times within $\pm$ 30 seconds of the associated calibration standard?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>XII. Target compound identification</b>				
Were relative retention times (RRT's) within $\pm$ 0.06 RRT units of the standard?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Did compound spectra meet specified EPA "Functional Guidelines" criteria?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were chromatogram peaks verified and accounted for?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>XIII. Compound quantitation/CRQLs</b>				
Were the correct internal standard (IS), quantitation ion and relative response factor (RRF) used to quantitate the compound?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were compound quantitation and CRQLs adjusted to reflect all sample dilutions and dry weight factors applicable to level I/IV validation?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>XIV. Tentatively identified compounds (TICs)</b>				
Were the major ions ( $> 10$ percent relative intensity) in the reference spectrum evaluated in sample spectrum?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	
Were relative intensities of the major ions within $\pm$ 20% between the sample and the reference spectra?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	
Did the raw data indicate that the laboratory performed a library search for all required peaks in the chromatograms (samples and blanks)?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	
<b>XV. System performance</b>				
System performance was found to be acceptable.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>XVI. Overall assessment of data</b>				
Overall assessment of data was found to be acceptable.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>XVII. Field duplicates</b>				
Field duplicate pairs were identified in this SDG.	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
Target compounds were detected in the field duplicates.	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	
<b>XVIII. Field blanks</b>				
Field blanks were identified in this SDG.	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
Target compounds were detected in the field blanks.	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	

**VALIDATION FINDINGS WORKSHEET**  
**Initial Calibration**

LDC #: 148-19C-1  
SDG #: 1PC-232X

**METHOD: GC/MS VOA (EPA SW 846 Method 8260B)**

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

N N/A Did the laboratory perform a 5 point calibration prior to sample analysis?

Y N N/A Were percent relative standard deviations (%RSD) and relative response factors (RRF) within method criteria for all CCC's and SPCC's?

Y N N/A Was a curve fit used for evaluation? If yes, what was the acceptance criteria used for evaluation? \_\_\_\_\_

Y N N/A Did the initial calibration meet the acceptance criteria?

Y N N/A Were all %RSDs and RRFs within the validation criteria of  $\leq 30\%$  RSD and  $\geq 0.05$  RRF?

#	Date	Standard ID	Compound	Finding %RSD (Limit: $\leq 30.0\%$ )	Finding RRF (Limit: $\geq 0.05$ )	Associated Samples	Qualifications
3/26/06	ICA-L		N		0.040	6003015-Bulk, 1	JWJ/A



LDC #: 14879C1  
 SDG #: 12e23y5

**VALIDATION FINDINGS WORKSHEET**  
**Blanks**

METHOD: GC/MS VOA (EPA SW 846 Method 8260B)

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

(Y) N N/A Was a method blank associated with every sample in this SDG?

(Y) N N/A Was a method blank analyzed at least once every 12 hours for each matrix and concentration?

(Y) N N/A Was there contamination in the method blanks? If yes, please see the qualifications below.

Blank analysis date: 4/3/06

Conc. units: ng/L

Associated Samples: 1

Compound	Blank ID	Sample Identification
Tetrahydrofuran	GPO3015-B1	1
Methylene chloride	3.42	6.1/20M
Acetone		
CRQL		

Blank analysis date: \_\_\_\_\_  
 Conc. units: \_\_\_\_\_

Associated Samples: \_\_\_\_\_

Compound	Blank ID	Sample Identification
Methylene chloride		
Acetone		
CRQL		

All results were qualified using the criteria stated below except those circled.

Note: Common contaminants such as Methylene chloride, Acetone, 2-Butanone, Carbon disulfide and TCs that were detected in samples within ten times the associated method blank concentration were qualified as not detected, "U". Other contaminants within five times the method blank concentration were also qualified as not detected, "U".

LDC #: 14X71C1  
SDG #: 1PC2375

**VALIDATION FINDINGS WORKSHEET**  
**Surrogate Spikes**

METHOD: GC/MS VOA (EPA SW 846 Method 8260B)

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".  
 N/N/A      Were all surrogate %R within QC limits?  
 N/N/A      If the percent recovery (%R) for one or more surrogates was out of QC limits, was a reanalysis performed to confirm samples with %R out of outside of criteria?

#	Date	Sample ID	Surrogate	%Recovery (Limits)	Qualifications
		1	BFB	(80-120)	↓ / <u>All except DV</u>
				( )	
				( )	
				( )	
				( )	
				( )	
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QC Limits (Soil)  
81-117  
88-110  
86-115  
80-120  
86-118

QC Limits (Water)  
74-121  
80-120  
80-120  
80-120

SMC1 (TOL) = Toluene-d8  
SMC2 (BFB) = Bromofluorobenzene  
SMC3 (DCE) = 1,2-Dichloroethane-d4  
SMC4 (DFM) = Dibromofluoromethane

LDC #: 1487A21  
SDG #: 1PC2325

**VALIDATION FINDINGS WORKSHEET**  
**Initial Calibration Calculation Verification**

Page: 1 of 1  
Reviewer: B  
2nd Reviewer: A

**METHOD: GC/MS VOA (EPA SW 846 Method 8260B)**

The Relative Response Factor (RRF), average RRF, and percent relative standard deviation (%RSD) were recalculated for the compounds identified below using the following calculations:

$$\text{RRF} = \frac{(A_x)(C_s)}{(A_b)(C_s)}$$

average RRF = sum of the RRFs/number of standards  
 $\%RSD = 100 * (\bar{X})$

$A_x$  = Area of compound,  
 $C_x$  = Concentration of compound,  
 $S$  = Standard deviation of the RRFs  
 $X$  = Mean of the RRFs

#	Standard ID	Calibration Date	Compound (Reference Internal Standard)	Reported	Recalculated	Reported	Recalculated	Reported	Recalculated
				RRF ( std)	RRF ( std)	Average RRF (Initial)	Average RRF (Initial)	%RSD	%RSD
1	3/26/06	1CAL	Methylene chloride (1st Internal standard)	0.484	0.484	0.498	0.498	5.05	5.05
			Trichlorethane (2nd Internal standard)	0.344	0.344	0.351	0.351	9.24	9.24
			Toluene (3rd Internal standard)	1.636	1.636	1.738	1.738	8.09	8.09
2			Methylene chloride (1st Internal standard)	1.471	1.471	1.481	1.481	10.98	10.98
			Trichlorethane (2nd Internal standard)						
			Toluene (3rd Internal standard)						
3			Methylene chloride (1st Internal standard)						
			Trichlorethane (2nd Internal standard)						
			Toluene (3rd Internal standard)						
4			Methylene chloride (1st Internal standard)						
			Trichlorethane (2nd Internal standard)						
			Toluene (3rd Internal standard)						

Comments: Refer to Initial Calibration findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

DC #: 148741  
DG #: 1PC2325

VALIDATION FINDINGS WORKSHEET  
Continuing Calibration Results Verification

Page: 1 of 1  
Reviewer: B  
2nd Reviewer: ef

METHOD: GC/MS VOA (EPA SW 846 Method 8260B)

The percent difference (%D) of the initial calibration average Relative Response Factors (RRFs) and the continuing calibration RRFs were recalculated for the compounds identified below using the following calculation:

$$\% \text{ Difference} = 100 * (\text{ave. RRF} - \text{RRF}) / \text{ave. RRF}$$
$$\text{RRF} = (A_x)(C_b) / (A_b)(C_x)$$

Where: ave. RRF = Initial calibration average RRF  
RRF = continuing calibration RRF

$A_x$  = Area of compound,

$C_x$  = Concentration of compound,

$A_b$  = Area of associated internal standard

$C_b$  = Concentration of internal standard

#	Standard ID	Calibration Date	Compound (Reference Internal Standard)	Reported		Recalculated	
				Average RRF (Initial)	RRF (CC)	RRF (CC)	%D
1	15T0025	4/1/06	<del>Chloro- methylene-chloride</del> (1st Internal standard)	25	20	23	12
			Trichlorethene (2nd Internal standard)	1.007	0.999	0.98	-0.8
			Toluene (3rd Internal standard)				
2	15T005A	4/3/06	Methylene chloride (1st Internal standard)	0.498	0.517	0.517	3.8
			Trichlorethene (2nd Internal standard)	0.351	0.353	0.353	0.6
			<del>Toluene</del> (3rd Internal standard)	1.738	1.622	1.622	6.7
3			Methylene chloride (1st Internal standard)	1.481	1.474	1.474	0.5
			Trichlorethene (2nd Internal standard)				
			Toluene (3rd Internal standard)				
4	15T008A	4/13/06	<del>Tetrachloro- methylene-chloride</del> (1st Internal standard)	0.063	0.065	0.065	3.2
			Trichlorethene (2nd Internal standard)				
			Toluene (3rd Internal standard)				

Comments: Refer to Continuing Calibration findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

LDC #: 1481PC1  
SDG #: 1PC2325

VALIDATION FINDINGS WORKSHEET  
Surrogate Results Verification

Page: \_\_\_\_/\_\_\_\_ of \_\_\_\_  
Reviewer: \_\_\_\_/\_\_\_\_  
2nd reviewer: \_\_\_\_/\_\_\_\_

METHOD: GC/MS VOA (EPA SW 846 Method 8260B)

The percent recoveries (%R) of surrogates were recalculated for the compounds identified below using the following calculation:

% Recovery: SF/SS \* 100

Where: SF = Surrogate Found  
SS = Surrogate Spiked

Sample ID: #1

	Surrogate Spiked	Surrogate Found	Percent Recovery Reported	Percent Recovery Recalculated	Percent Difference
Toluene-d8	25	25.17	101	101	0
Bromofluorobenzene	25	31.99	128	128	0
1,2-Dichloroethane-d4					
Dibromofluoromethane	25	27.83	111	111	0

Sample ID:

	Surrogate Spiked	Surrogate Found	Percent Recovery Reported	Percent Recovery Recalculated	Percent Difference
Toluene-d8					
Bromofluorobenzene					
1,2-Dichloroethane-d4					
Dibromofluoromethane					

Sample ID:

	Surrogate Spiked	Surrogate Found	Percent Recovery Reported	Percent Recovery Recalculated	Percent Difference
Toluene-d8					
Bromofluorobenzene					
1,2-Dichloroethane-d4					
Dibromofluoromethane					

Sample ID:

	Surrogate Spiked	Surrogate Found	Percent Recovery Reported	Percent Recovery Recalculated	Percent Difference
Toluene-d8					
Bromofluorobenzene					
1,2-Dichloroethane-d4					
Dibromofluoromethane					

Sample ID:

	Surrogate Spiked	Surrogate Found	Percent Recovery Reported	Percent Recovery Recalculated	Percent Difference
Toluene-d8					
Bromofluorobenzene					
1,2-Dichloroethane-d4					
Dibromofluoromethane					

LDC #: 14819C  
SDG #: 19C2325

## VALIDATION FINDINGS WORKSHEET Laboratory Control Sample Results Verification

### METHOD: GC/MS VOA (EPA SW 846 Method 8260B)

The percent recoveries (%R) and Relative Percent Difference (RPD) of the laboratory control sample and laboratory control sample duplicate (if applicable) were recalculated for the compounds identified below using the following calculation:

$$\% \text{ Recovery} = 100 * \text{SSC/SA}$$

Where:  
SSC = Spiked sample concentration  
SA = Spike added

$$\text{RPD} = | \text{LCS} - \text{LCSD} | * 2 / (\text{LCS} + \text{LCSD})$$

LCS ID: 6903015-P51

LCS = Laboratory control sample percent recovery

LCSD = Laboratory control sample duplicate percent recovery

Compound	Spiked Added ( <u>ug/L</u> )	Spiked Sample Concentration ( <u>ug/L</u> )		Percent Recovery	LCS	LCSD	Percent Recovery		RPD		LCS/LCSD	Comments:
		LCS	LCSD		Reported	Recalc.	Reported	Recalc.	Reported	Recalc.		
1,1-Dichloroethene	25.0	NA	25.1	NA	100	100	100	100	0.0	0.0	1.00	Refer to Laboratory Control Sample findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.
Trichloroethene			26.0		100	100	100	100				
Benzene			24.2		97	97	97	97				
Toluene			23.7		95	95	95	95				
Chlorobenzene			24.5		98	98	98	98				

Comments: Refer to Laboratory Control Sample findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

LDC #: 14819C  
 SDG #: 1PC 2325

**VALIDATION FINDINGS WORKSHEET**  
Sample Calculation Verification

Page: 1 of 1

Reviewer: JF

2nd reviewer: AL

**METHOD:** GC/MS VOA (EPA SW 846 Method 8260B)

Were all reported results recalculated and verified for all level IV samples?

Were all recalculated results for detected target compounds agree within 10.0% of the reported results?

$$\text{Concentration} = \frac{(A_s)(L)(DF)}{(A_u)(RRF)(V_o)(\%)S}$$

$A_u$  = Area of the characteristic ion (EICP) for the compound to be measured

$A_s$  = Area of the characteristic ion (EICP) for the specific internal standard

$I_s$  = Amount of internal standard added in nanograms (ng)

RRF = Relative response factor of the calibration standard.

$V_o$  = Volume or weight of sample pruged in milliliters (ml) or grams (g).

Df = Dilution factor.

$\%S$  = Percent solids, applicable to soils and solid matrices only.

Example:

Sample I.D. #1, 1, u = Dichlorobenzene

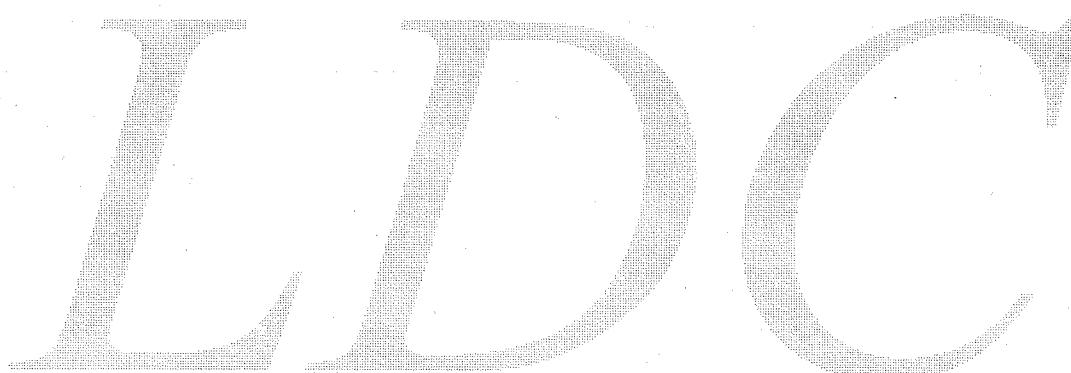
$$\text{Conc.} = \frac{(147977)(25)}{(509472)(1.639)} =$$

8.9 ug/L

#	Sample ID	Compound	Reported Concentration ( )	Calculated Concentration ( )	Qualification

**Boeing Realty Corp., Former C-6 Facility, Torrance CA**  
**Data Validation Reports**  
**LDC# 14879**

**Manganese**



## **Laboratory Data Consultants, Inc.**

### **Data Validation Report**

**Project/Site Name:** Boeing Realty Corp, Former C-6 Facility, Torrance CA

**Collection Date:** March 20, 2006

**LDC Report Date:** May 15, 2006

**Matrix:** Water

**Parameters:** Manganese

**Validation Level:** Tier 1

**Laboratory:** Del Mar Analytical

**Sample Delivery Group (SDG):** IPC2080

#### **Sample Identification**

MWC011\_WG032006\_0001

## **Introduction**

This data review covers one water sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 6010B for Manganese.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) as there are no current guidelines for the methods stated above.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified a P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blanks are summarized in Section III.

Field duplicates are summarized in Section XIII.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- U     Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
  - J     Indicates an estimated value.
  - R     Quality control indicates the data is not usable.
  - N     Presumptive evidence of presence of the constituent.
  - UJ    Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
  - A     Indicates the finding is based upon technical validation criteria.
  - P     Indicates the finding is related to a protocol/contractual deviation.
- None    Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

Calibration data were not reviewed for Tier 1.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the initial, continuing and preparation blanks.

## **IV. ICP Interference Check Sample (ICS) Analysis**

The frequency of analysis was met.

The criteria for analysis were met.

## **V. Matrix Spike Analysis**

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VI. Duplicate Sample Analysis**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable.

## **VII. Laboratory Control Samples (LCS)**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits.

## **VIII. Internal Standards (ICP-MS)**

ICP-MS was not utilized in this SDG.

## **IX. Furnace Atomic Absorption QC**

Graphite furnace atomic absorption was not utilized in this SDG.

## **X. ICP Serial Dilution**

ICP serial dilution was not performed for this SDG.

## **XI. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **XII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **XIII. Field Duplicates**

No field duplicates were identified in this SDG.

## **XIV. Field Blanks**

No field blanks were identified in this SDG.

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Manganese - Data Qualification Summary - SDG IPC2080**

No Sample Data Qualified in this SDG

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Manganese - Laboratory Blank Data Qualification Summary - SDG IPC2080**

No Sample Data Qualified in this SDG



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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2080

Sampled: 03/20/06  
 Received: 03/20/06

### DISSOLVED METALS

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
<b>Sample ID: IPC2080-07 (MWC011_WG032006_0001 - Water)</b>									
Reporting Units: mg/l									
Manganese	EPA 6010B-Diss	6C20131	0.0070	0.020	0.25	1	03/20/06	03/23/06	
<b>Sample ID: IPC2080-11 (MWB019_WG032006_0001 - Water)</b>									
Reporting Units: mg/l									
Manganese	EPA 6010B-Diss	6C20131	0.0070	0.020	0.0094	1	03/20/06	03/23/06	J

**Del Mar Analytical - Irvine**  
 Michele Chamberlin  
 Project Manager

ML 251506

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LDC #: 14879A4

SDG #: IPC2080

Laboratory: Del Mar Analytical

**VALIDATION COMPLETENESS WORKSHEET**

Tier 1

Date: 5/15/06

Page: 1 of 1

Reviewer: MU

2nd Reviewer: A

**METHOD:** Manganese (EPA SW 846 Method 6010B)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	<b>Validation Area</b>		<b>Comments</b>
I.	Technical holding times	A	Sampling dates: 3/20/06
II.	Calibration	N	
III.	Blanks	A	
IV.	ICP Interference Check Sample (ICS) Analysis	N	
V.	Matrix Spike Analysis	A	yes/no in client sample
VI.	Duplicate Sample Analysis	N	
VII.	Laboratory Control Samples (LCS)	A	LCS
VIII.	Internal Standard (ICP-MS)	N	Not utilized
IX.	Furnace Atomic Absorption QC	N	
X.	ICP Serial Dilution	N	Not performed
XI.	Sample Result Verification	N	
XII.	Overall Assessment of Data	A	
XIII.	Field Duplicates	N	
XIV.	Field Blanks	N	

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet

ND = No compounds detected  
 R = Rinsate  
 FB = Field blank

D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples: 182

1	MWC011_WG032006_0001	11		21		31	
2	RB	12		22		32	
3		13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

## **Laboratory Data Consultants, Inc.**

### **Data Validation Report**

**Project/Site Name:** Boeing Realty Corp, Former C-6 Facility, Torrance CA

**Collection Date:** March 21, 2006

**LDC Report Date:** May 15, 2006

**Matrix:** Water

**Parameters:** Manganese

**Validation Level:** Tier 2

**Laboratory:** Del Mar Analytical

**Sample Delivery Group (SDG):** IPC2190

#### **Sample Identification**

IRZB0095\_WG032106\_0001

## **Introduction**

This data review covers one water sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 6010B for Manganese.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) as there are no current guidelines for the methods stated above.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified a P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blanks are summarized in Section III.

Field duplicates are summarized in Section XIII.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- U     Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
  - J     Indicates an estimated value.
  - R     Quality control indicates the data is not usable.
  - N     Presumptive evidence of presence of the constituent.
  - UJ    Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
  - A     Indicates the finding is based upon technical validation criteria.
  - P     Indicates the finding is related to a protocol/contractual deviation.
- None    Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

An initial calibration was performed.

The frequency and analysis criteria of the initial calibration verification (ICV) and continuing calibration verification (CCV) were met.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the initial, continuing and preparation blanks.

## **IV. ICP Interference Check Sample (ICS) Analysis**

The frequency of analysis was met.

The criteria for analysis were met.

## **V. Matrix Spike Analysis**

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VI. Duplicate Sample Analysis**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable.

## **VII. Laboratory Control Samples (LCS)**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits.

## **VIII. Internal Standards (ICP-MS)**

ICP-MS was not utilized in this SDG.

## **IX. Furnace Atomic Absorption QC**

Graphite furnace atomic absorption was not utilized in this SDG.

## **X. ICP Serial Dilution**

ICP serial dilution was not performed for this SDG.

## **XI. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **XII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **XIII. Field Duplicates**

No field duplicates were identified in this SDG.

## **XIV. Field Blanks**

No field blanks were identified in this SDG.

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Manganese - Data Qualification Summary - SDG IPC2190**

No Sample Data Qualified in this SDG

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Manganese - Laboratory Blank Data Qualification Summary - SDG IPC2190**

No Sample Data Qualified in this SDG



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 2520 E. Sunset Rd. #3, Las Vegas, NV 89120 (702) 798-3620 FAX (702) 798-3621

TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2190

Sampled: 03/21/06  
 Received: 03/21/06

### DISSOLVED METALS

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
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**Sample ID: IPC2190-03 (IRZMW001B\_WG032106\_0001 - Water)**

Reporting Units: mg/l

Manganese	EPA 6010B-Diss	6C22121	0.0070	0.020	0.025	1	03/22/06	03/23/06
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**Sample ID: IPC2190-04 (IRZB0095\_WG032106\_0001 - Water)**

Reporting Units: mg/l

Manganese	EPA 6010B-Diss	6C22121	0.0070	0.020	1.7	1	03/22/06	03/23/06
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**Sample ID: IPC2190-09 (WCC\_6S\_WG032106\_0001 - Water)**

Reporting Units: mg/l

Manganese	EPA 6010B-Diss	6C22121	0.0070	0.020	6.3	1	03/22/06	03/23/06
-----------	----------------	---------	--------	-------	-----	---	----------	----------

**Sample ID: IPC2190-11 (CMW026\_WG032106\_0001 - Water)**

Reporting Units: mg/l

Manganese	EPA 6010B-Diss	6C22121	0.0070	0.020	0.064	1	03/22/06	03/23/06
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**Del Mar Analytical - Irvine**  
 Michele Chamberlin  
 Project Manager

16-051596

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IPC2190 <Page 29 of 56>

LDC #: 14879B4  
SDG #: IPC2190  
Laboratory: Del Mar Analytical

## VALIDATION COMPLETENESS WORKSHEET

Tier 2

Date: 5/15/06

Page: 1 of 1

Reviewer: MM

2nd Reviewer: A

METHOD: Manganese (EPA SW 846 Method 6010B)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 3/11/06
II.	Calibration	A	
III.	Blanks	A	
IV.	ICP Interference Check Sample (ICS) Analysis	A	
V.	Matrix Spike Analysis	A	MS/MSD ICP-2190-3
VI.	Duplicate Sample Analysis	N	
VII.	Laboratory Control Samples (LCS)	A	LCS
VIII.	Internal Standard (ICP-MS)	N	34 wt % Mg
IX.	Furnace Atomic Absorption QC	N	
X.	ICP Serial Dilution	N	Not performed.
XI.	Sample Result Verification	N	
XII.	Overall Assessment of Data	A	
XIII.	Field Duplicates	U	
XIV.	Field Blanks	N	

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:

1	IRZB0095_WG032106_0001	11		21		31	
2	PB	12		22		32	
3		13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

## **Laboratory Data Consultants, Inc.**

### **Data Validation Report**

**Project/Site Name:** Boeing Realty Corp, Former C-6 Facility, Torrance CA

**Collection Date:** March 22, 2006

**LDC Report Date:** May 15, 2006

**Matrix:** Water

**Parameters:** Manganese

**Validation Level:** Tier 3

**Laboratory:** Del Mar Analytical

**Sample Delivery Group (SDG):** IPC2325

#### **Sample Identification**

CMW0002\_WG032206\_0001

## **Introduction**

This data review covers one water sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Method 6010B for Manganese.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) as there are no current guidelines for the methods stated above.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified a P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blanks are summarized in Section III.

Field duplicates are summarized in Section XIII.

The following are definitions of the data qualifiers:

- U     Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- J     Indicates an estimated value.
- R     Quality control indicates the data is not usable.
- N     Presumptive evidence of presence of the constituent.
- UJ    Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A     Indicates the finding is based upon technical validation criteria.
- P     Indicates the finding is related to a protocol/contractual deviation.

None    Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

An initial calibration was performed.

The frequency and analysis criteria of the initial calibration verification (ICV) and continuing calibration verification (CCV) were met.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the initial, continuing and preparation blanks.

## **IV. ICP Interference Check Sample (ICS) Analysis**

The frequency of analysis was met.

The criteria for analysis were met.

## **V. Matrix Spike Analysis**

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **VI. Duplicate Sample Analysis**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable.

## **VII. Laboratory Control Samples (LCS)**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits.

## **VIII. Internal Standards (ICP-MS)**

ICP-MS was not utilized in this SDG.

## **IX. Furnace Atomic Absorption QC**

Graphite furnace atomic absorption was not utilized in this SDG.

## **X. ICP Serial Dilution**

ICP serial dilution was not performed for this SDG.

## **XI. Sample Result Verification**

All sample result verifications were acceptable.

## **XII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **XIII. Field Duplicates**

No field duplicates were identified in this SDG.

## **XIV. Field Blanks**

No field blanks were identified in this SDG.

**Boeing Realty Corp, Former C-6 Facility, Torrance CA  
Manganese - Data Qualification Summary - SDG IPC2325**

No Sample Data Qualified in this SDG

**Boeing Realty Corp, Former C-6 Facility, Torrance CA  
Manganese - Laboratory Blank Data Qualification Summary - SDG IPC2325**

No Sample Data Qualified in this SDG



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2520 E. Sunset Rd. #3, Las Vegas, NV 89120 (702) 798-3620 FAX (702) 798-3621

TAIT Environmental/Boeing  
701 N. Parkcenter Drive  
Santa Ana, CA 92705  
Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
EM2303D/PHASE 01  
Report Number: IPC2325

Sampled: 03/22/06  
Received: 03/22/06

### DISSOLVED METALS

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Date Qualifiers
<b>Sample ID: IPC2325-04 (IRZCMW002_WG032206_0001 - Water)</b>									
Reporting Units: mg/l									
Manganese	EPA 6010B-Diss	6C23108	0.0070	0.020	3.8	1	03/23/06	03/23/06	
<b>Sample ID: IPC2325-09 (CMW0002_WG032206_0001 - Water)</b>									
Reporting Units: mg/l									
Manganese	EPA 6010B-Diss	6C23108	0.0070	0.020	0.21	1	03/23/06	03/23/06	

Del Mar Analytical - Irvine  
Michele Chamberlin  
Project Manager

LOREK/JD6

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IPC2325 <Page 30 of 73>

LDC #: 14879C4

SDG #: IPC2325

Laboratory: Del Mar Analytical

**VALIDATION COMPLETENESS WORKSHEET**

Tier 3

Date: 5/15/06

Page: 1 of 1

Reviewer: J.W.

2nd Reviewer: R.

**METHOD:** Manganese (EPA SW 846 Method 6010B)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	<b>Validation Area</b>		<b>Comments</b>
I.	Technical holding times	A	Sampling dates: 3/22/06
II.	Calibration	A	
III.	Blanks	A	
IV.	ICP Interference Check Sample (ICS) Analysis	A	
V.	Matrix Spike Analysis	A	MS/MS No client sample.
VI.	Duplicate Sample Analysis	N	
VII.	Laboratory Control Samples (LCS)	A	LCS
VIII.	Internal Standard (ICP-MS)	N	3 wt % Iridium
IX.	Furnace Atomic Absorption QC	N	
X.	ICP Serial Dilution	N	Not prepared
XI.	Sample Result Verification	A	
XII.	Overall Assessment of Data	A	
XIII.	Field Duplicates	N	
XIV.	Field Blanks	N	

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet

ND = No compounds detected  
 R = Rinsate  
 FB = Field blank

D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples: No

1	CMW0002_WG032206_0001	11		21		31	
2	PR	12		22		32	
3		13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

LDC #: 14879c4  
SDG #: 1PL2325

VALIDATION FINDINGS CHECKLIST

Page: 1 of 2  
Reviewer: LM  
2nd Reviewer: R

**Method:** Metals (EPA SW 846 Method 6010/7000/6020)

Validation Area	Yes	No	NA	Findings/Comments
<b>I. Technical holding times</b>				
All technical holding times were met.	✓			
Cooler temperature criteria was met.	✓			
<b>II. Calibration</b>				
Were all instruments calibrated daily, each set-up time?	✓			
Were the proper number of standards used?	✓			
Were all initial and continuing calibration verification %Rs within the 90-110% (80-120% for mercury and 85-115% for cyanide) QC limits?	✓			
Were all initial calibration correlation coefficients > 0.995? (Level IV only)	✓			
<b>III. Blanks</b>				
Was a method blank associated with every sample in this SDG?	✓			
Was there contamination in the method blanks? If yes, please see the Blanks validation completeness worksheet.		✓		
<b>IV. ICP Interference Check Sample</b>				
Were ICP interference check samples performed daily?	✓			
Were the AB solution percent recoveries (%R) with the 80-120% QC limits?	✓			
<b>V. Matrix spike/Matrix spike duplicates</b>				
Were a matrix spike (MS) and duplicate (DUP) analyzed for each matrix in this SDG? If no, indicate which matrix does not have an associated MS/MSD or MS/DUP. Soil / Water.	✓			
Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the 75-125 QC limits? If the sample concentration exceeded the spike concentration by a factor of 4 or more, no action was taken.	✓			
Were the MS/MSD or duplicate relative percent differences (RPD) ≤ 20% for waters and ≤ 35% for soil samples? A control limit of +/- RL (+/- 2X RL for soil) was used for samples that were ≤ 5X the RL, including when only one of the duplicate sample values were < 5X the RL.	✓			
<b>VI. Laboratory control samples</b>				
Was an LCS analyzed for this SDG?	✓			
Was an LCS analyzed per extraction batch?	✓			
Were the LCS percent recoveries (%R) and relative percent difference (RPD) within the 80-120% QC limits for water samples and laboratory established QC limits for soils?	✓			
<b>VII. Furnace Atomic Absorption QC</b>				
If MSA was performed, was the correlation coefficients > 0.995?			✓	
Do all applicable analyses have duplicate injections? (Level IV only)			✓	
For sample concentrations > RL, are applicable duplicate injection RSD values < 20%? (Level IV only)			✓	
Were analytical spike recoveries within the 85-115% QC limits?			✓	

LDC #: 14819 C4  
SDG #: 7D-2325

VALIDATION FINDINGS CHECKLIST

Page: 2 of ✓  
Reviewer: JM  
2nd Reviewer: C

Validation Area	Yes	No	NA	Findings/Comments
<b>VII. ICP Serial Dilution</b>				
Was an ICP serial dilution analyzed if analyte concentrations were > 50X the IDL?	✓	X		
Were all percent differences (%Ds) < 10%?		✓		
Was there evidence of negative interference? If yes, professional judgement will be used to qualify the data.		✓		
<b>VIII. Internal Standards (EPA SW 846 Method 6020)</b>				
Were all the percent recoveries (%R) within the 30-120% of the intensity of the internal standard in the associated initial calibration?		✓		
If the %Rs were outside the criteria, was a reanalysis performed?		✓		
<b>IX. Regional Quality Assurance and Quality Control</b>				
Were performance evaluation (PE) samples performed?		✓		
Were the performance evaluation (PE) samples within the acceptance limits?		✓		
<b>X. Sample Result Verification</b>				
Were RLs adjusted to reflect all sample dilutions and dry weight factors applicable to level IV validation?	✓			
<b>XI. Overall assessment of data</b>				
Overall assessment of data was found to be acceptable.	✓			
<b>XII. Field duplicates</b>				
Field duplicate pairs were identified in this SDG.		✓		
Target analytes were detected in the field duplicates.				
<b>XIII. Field blanks</b>				
Field blanks were identified in this SDG.		✓		
Target analytes were detected in the field blanks.		✓		

LDC #: 14819cc4  
SDG #: TPC225

**VALIDATION FINDINGS WORKSHEET**  
**Initial and Continuing Calibration Calculation Verification**

Page: 1 of 1  
Reviewer: W  
2nd Reviewer: K

**METHOD:** Trace Metals (EPA SW 846 Method 6010/7000)

An initial and continuing calibration verification percent recovery (%R) was recalculated for each type of analysis using the following formula:

$$\%R = \frac{\text{Found}}{\text{True}} \times 100$$

Where, Found = concentration (in ug/L) of each analyte measured in the analysis of the ICV or CCV solution  
True = concentration (in ug/L) of each analyte in the ICV or CCV source

Standard ID	Type of Analysis	Element	Found (ug/L)	True (ug/L)	Recalculated %R	Reported %R	Acceptable (Y/N)
Tn	ICP (Initial calibration)	Mn	2.014	2.00	101	101	Y
	GFAA (Initial calibration)						
	CVAA (Initial calibration)						
ccv	ICP (Continuing calibration)	Mn	1.001	1.0	100	100	Y
	GFAA (Continuing calibration)						
	CVAA (Continuing calibration)						
	Cyanide (Initial calibration)						
	Cyanide (Continuing calibration)						

Comments: Refer to Calibration Verification findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

LDC #: 148194  
SDG #: 7PL2325

**VALIDATION FINDINGS WORKSHEET**  
**Level IV Recalculation Worksheet**

Page: 1 of 1  
Reviewer: MW  
2nd Reviewer:  

**METHOD:** Trace Metals (EPA SW 846 Method 6010/7000)

Percent recoveries (%R) for an ICP interference check sample, a laboratory control sample and a matrix spike sample were recalculated using the following formula:

$$\%R = \frac{\text{Found}}{\text{True}} \times 100$$
 Where, Found = Concentration of each analyte measured in the analysis of the sample. For the matrix spike calculation,  

$$True = \frac{\text{SSR} - \text{SR}}{\text{SSR}}$$
 Found = SSR (spiked sample result) - SR (sample result).  
 True = Concentration of each analyte in the source.

A sample and duplicate relative percent difference (RPD) was recalculated using the following formula:

$$RPD = \frac{|S-D|}{(S+D)/2} \times 100$$
 Where, S = Original sample concentration  
 D = Duplicate sample concentration

An ICP serial dilution percent difference (%D) was recalculated using the following formula:

$$\%D = \frac{|I-SDR|}{SDR} \times 100$$
 Where, I = Initial Sample Result (mg/l)  
 SDR = Serial Dilution Result (mg/l) (Instrument Reading  $\times 5$ )

Sample ID	Type of Analysis	Element	Found / S / I (units)	True / D / SDR (units)	Reported		Acceptable (Y/N)
					Recalculated	%R / RPD / %D	
Trace	ICP Interference check	Hg	0.4681	0.500	94	RP	Y
LCS	Laboratory control sample		0.9798	1.00	98	98	
7PL2325	Matrix spike	(SSR-SR)	0.9367	1	94	94	
7PL2325	Duplicate		0.9157	0.9367	2	2	
MX	ICP serial dilution						

Comments: Refer to appropriate worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

LDC #: 14879 c4  
SDG #: JPC2325

## **VALIDATION FINDINGS WORKSHEET**

### **Sample Calculation Verification**

Page: 1 of 1  
Reviewer: MV  
2nd reviewer: N

**METHOD: Trace Metals (EPA SW 846 Method 6010/7000)**

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

- N N/A Have results been reported and calculated correctly?  
 N N/A Are results within the calibrated range of the instruments and within the linear range of the ICP?  
 N N/A Are all detection limits below the CRDL?

Detected analyte results for \_\_\_\_\_ were recalculated and verified using the following equation:

**Concentration =**  $\frac{(RD)(FV)(Dil)}{(In. Vol.)(\%)S}$  **Recalculation:**

#### **Recalculation:**

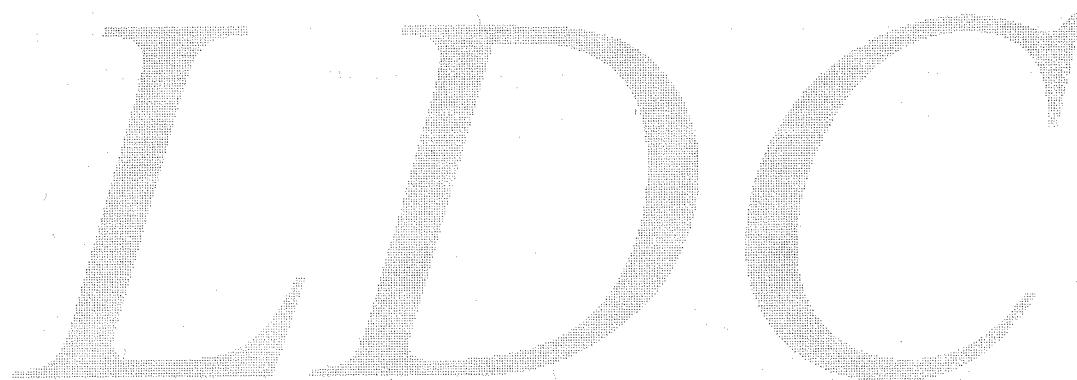
RD	=	Raw data concentration
FV	=	Final volume (ml)
In. Vol.	=	Initial volume (ml) or weight (G)
Dil	=	Dilution factor
%S	=	Decimal percent solids

From the new lot

$$M_4 = 0.2057 \text{ mg/L}$$

**Boeing Realty Corp., Former C-6 Facility, Torrance CA**  
**Data Validation Reports**  
**LDC# 14879**

**Wet Chemistry**



## **Laboratory Data Consultants, Inc.**

### **Data Validation Report**

**Project/Site Name:** Boeing Realty Corp, Former C-6 Facility, Torrance CA

**Collection Date:** March 20, 2006

**LDC Report Date:** May 15, 2006

**Matrix:** Water

**Parameters:** Wet Chemistry

**Validation Level:** Tier 1

**Laboratory:** Del Mar Analytical

**Sample Delivery Group (SDG):** IPC2080

#### **Sample Identification**

MWC011\_WG032006\_0001  
IRZMW002B\_WG032006\_0001  
IRZMW002B\_WG032006\_0001DUP

## **Introduction**

This data review covers 3 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 300.0 for Chloride, Nitrate, Nitrite, Orthophosphate, and Sulfate, EPA Method 310.1 for Alkalinity, EPA Method 350.3 for Ammonia, EPA Method 415.1 for Total Organic Carbon, and Standard Method 4500-CO<sub>2</sub> for Carbon Dioxide.

The review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) as there are no current guidelines for the methods stated above.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified a P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blank results are summarized in Section III.

Field duplicates are summarized in Section IX.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- U     Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- J     Indicates an estimated value.
- R     Quality control indicates the data is not usable.
- N     Presumptive evidence of presence of the constituent.
- UJ    Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A     Indicates the finding is based upon technical validation criteria.
- P     Indicates the finding is related to a protocol/contractual deviation.

None    Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodices were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration data were not reviewed for Tier I.

### **b. Calibration Verification**

Calibration verification data were not reviewed for Tier I.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the method blanks with the following exceptions:

Method Blank ID	Analyte	Concentration	Associated Samples
MB	Chloride	0.164 mg/L	MWC011_WG032006_0001

Sample concentrations were compared to concentrations detected in the method blanks. The sample concentrations were either not detected or were significantly greater ( $>5X$  blank contaminants) than the concentrations found in the associated method blanks.

## **IV. Matrix Spike/Matrix Spike Duplicates**

Matrix spike (MS) and matrix spike duplicate (MSD) analyses were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **V. Duplicates**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable. Relative percent differences (RPD) were within QC limits.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits.

## **VII. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

## **X. Field Blanks**

No field blanks were identified in this SDG.

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Wet Chemistry - Data Qualification Summary - SDG IPC2080**

No Sample Data Qualified in this SDG

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG IPC2080**

No Sample Data Qualified in this SDG



**Del Mar Analytical**

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 1014 E. Cooley Dr., Suite A, Colton, CA 92324 (909) 370-4667 FAX (909) 370-1046  
 9830 South 51st St., Suite B-120, Phoenix, AZ 85044 (480) 785-0043 FAX (480) 785-0851  
 2520 E. Sunset Rd. #3, Las Vegas, NV 89120 (702) 798-3620 FAX (702) 798-3621

TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2080

Sampled: 03/20/06  
 Received: 03/20/06

## INORGANICS

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
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**Sample ID: IPC2080-07 (MWC011\_WG032006\_0001 - Water)**

Reporting Units: mg/l

Carbon Dioxide	SM4500-CO2	6C20139	2.0	2.0	5.3	1	03/20/06	03/20/06	
Ammonia-NH3	EPA 350.3	6C27056	0.14	0.60	ND	1	03/24/06	03/24/06	
Alkalinity as CaCO3	EPA 310.1	6C23139	2.0	2.0	180	1	03/23/06	03/23/06	
Chloride	EPA 300.0	6C20055	1.5	5.0	200	10	03/20/06	03/20/06	
Nitrate-NO3	EPA 300.0	6C20055	0.35	0.50	5.2	1	03/20/06	03/20/06	
Nitrite-NO2	EPA 300.0	6C20055	3.0	5.0	ND	10	03/20/06	03/20/06	RL-1
Orthophosphate - PO4	EPA 300.0	6C20055	0.40	0.50	ND	1	03/20/06	03/20/06	
Sulfate	EPA 300.0	6C20055	0.45	0.50	20	1	03/20/06	03/20/06	
Total Organic Carbon	EPA 415.1	6C23066	0.25	1.0	5.3	1	03/22/06	03/22/06	

**Sample ID: IPC2080-08 (IRZMW002B\_WG032006\_0001 - Water)**

Reporting Units: mg/l

Sulfide	EPA 376.2	6C21070	0.010	0.10	ND	1	03/21/06	03/21/06
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**Sample ID: IPC2080-09 (IRZMW002B\_WG032006\_0002 - Water)**

Reporting Units: mg/l

Sulfide	EPA 376.2	6C21070	0.010	0.10	ND	1	03/21/06	03/21/06
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**Sample ID: IPC2080-10 (IRZB0081\_WG032006\_0001 - Water)**

Reporting Units: mg/l

Sulfide	EPA 376.2	6C21070	0.010	0.10	ND	1	03/21/06	03/21/06
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**Sample ID: IPC2080-11 (MWB019\_WG032006\_0001 - Water)**

Reporting Units: mg/l

Carbon Dioxide	SM4500-CO2	6C20139	2.0	2.0	62	1	03/20/06	03/20/06	
Ammonia-NH3	EPA 350.3	6C27056	0.14	0.60	ND	1	03/24/06	03/24/06	
Alkalinity as CaCO3	EPA 310.1	6C23139	2.0	2.0	390	1	03/23/06	03/23/06	
Chloride	EPA 300.0	6C20055	3.0	10	260	20	03/20/06	03/20/06	
Nitrate-NO3	EPA 300.0	6C20055	7.0	10	90	20	03/20/06	03/20/06	
Nitrite-NO2	EPA 300.0	6C20055	6.0	10	ND	20	03/20/06	03/20/06	RL-1
Orthophosphate - PO4	EPA 300.0	6C20055	0.40	0.50	ND	1	03/20/06	03/20/06	
Sulfate	EPA 300.0	6C20055	9.0	10	610	20	03/20/06	03/20/06	
Total Organic Carbon	EPA 415.1	6C23066	0.25	1.0	2.0	1	03/22/06	03/22/06	

**Del Mar Analytical - Irvine**  
 Michele Chamberlin  
 Project Manager

*M. Chamberlin*

The results pertain only to the samples tested in the laboratory. This report shall not be reproduced,  
 except in full, without written permission from Del Mar Analytical.

IPC2080 <Page 28 of 73>

BOE-C6-0050765

LDC #: 14879A6

## VALIDATION COMPLETENESS WORKSHEET

Date: 6/5/06

SDG #: IPC2080

Tier 1

Page: 1 of 1

Laboratory: Del Mar Analytical

Reviewer: MY

2nd Reviewer: 

~~o POC~~ ~~POC~~ ~~143~~  
**METHOD: (Analyte)** Alkalinity (EPA Method 310.1), Ammonia-N (EPA Method 350.3), Chloride, Nitrate-X, Nitrite-X, Orthophosphate-X, Sulfate (EPA Method 300.0), TOC (EPA Method 415.1), Carbon Dioxide (SM4500+ CO<sub>2</sub>)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 3/20/06
IIa.	Initial calibration	N	
IIb.	Calibration verification	N	
III.	Blanks	SW	
IV	Matrix Spike/Matrix Spike Duplicates	A	3 Ls Musb / dup.
V	Duplicates	A	
VI.	Laboratory control samples	A	Ls
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N <del>ND</del> A	(2, IRZMW002B-WG032006-0002)
X	Field blanks	N	

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet

ND = No compounds detected  
 R = Rinsate  
 FB = Field blank

D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples:

1	MWC011_WG032006_0001	11		21		31	
2	IRZMW002B_WG032006_0001	12		22		32	
3	↓ PUP	13		23		33	
4	14B	14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes: \_\_\_\_\_  
 \_\_\_\_\_

LDC #: 14879A6  
SDG #: TPLC2080

**VALIDATION FINDINGS WORKSHEET**  
**Sample Specific Analysis Reference**

Page: 1 of 1  
Reviewer: WY  
2nd reviewer: A

All circled methods are applicable to each sample.

Sample ID	Parameter
1	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> <u>Cl<sub>2</sub></u>
2	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup> <u>S</u>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>6+</sup>

Comments:

LDC #: 1487946  
SDG #: TPLC2080

# VALIDATION FINDINGS WORKSHEET

Blanks

## METHOD: Inorganics, Me

See over

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

N/A Were all samples associated with a given method blank?

Conc. units: mg/l

### **Associated Samples:**

1 (75x)

Page: 1 of 1  
Reviewer: luny  
2nd Reviewer: lc

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:  
All contaminants within five times the method blank concentration were qualified as not detected. "u."

**Laboratory Data Consultants, Inc.**  
**Data Validation Report**

**Project/Site Name:** Boeing Realty Corp, Former C-6 Facility, Torrance CA  
**Collection Date:** March 21, 2006  
**LDC Report Date:** May 15, 2006  
**Matrix:** Water  
**Parameters:** Wet Chemistry  
**Validation Level:** Tier 2  
**Laboratory:** Del Mar Analytical

**Sample Delivery Group (SDG):** IPC2190

**Sample Identification**

IRZB0095\_WG032106\_0001

## Introduction

This data review covers one water sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 300.0 for Chloride, Nitrate, Nitrite, Orthophosphate, and Sulfate, EPA Method 310.1 for Alkalinity, EPA Method 350.3 for Ammonia, EPA Method 376.2 for Sulfide, EPA Method 415.1 for Total Organic Carbon, and Standard Method 4500-CO<sub>2</sub> for Carbon Dioxide.

The review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) as there are no current guidelines for the methods stated above.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified a P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blank results are summarized in Section III.

Field duplicates are summarized in Section IX.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- U     Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- J     Indicates an estimated value.
- R     Quality control indicates the data is not usable.
- N     Presumptive evidence of presence of the constituent.
- UJ    Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A     Indicates the finding is based upon technical validation criteria.
- P     Indicates the finding is related to a protocol/contractual deviation.
- None   Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodices were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

All criteria for the initial calibration of each method were met.

### **b. Calibration Verification**

Calibration verification frequency and analysis criteria were met for each method when applicable.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the method blanks with the following exceptions:

Method Blank ID	Analyte	Concentration	Associated Samples
ICB/CCB	Chloride	0.18 mg/L	All samples in SDG IPC2190

Sample concentrations were compared to concentrations detected in the method blanks. The sample concentrations were either not detected or were significantly greater ( $>5X$  blank contaminants) than the concentrations found in the associated method blanks.

## **IV. Matrix Spike/Matrix Spike Duplicates**

Matrix spike (MS) and matrix spike duplicate (MSD) analyses were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **V. Duplicates**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable. Relative percent differences (RPD) were within QC limits.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits.

## **VII. Sample Result Verification**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

## **X. Field Blanks**

No field blanks were identified in this SDG.

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Wet Chemistry - Data Qualification Summary - SDG IPC2190**

No Sample Data Qualified in this SDG

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG IPC2190**

No Sample Data Qualified in this SDG



**Del Mar Analytical**

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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2190

Sampled: 03/21/06  
 Received: 03/21/06

## INORGANICS

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
<b>Sample ID: IPC2190-03 (HRZMW001B_WG032106_0001 - Water)</b>									
Reporting Units: mg/l									
Carbon Dioxide	SM4500-CO2	6C21139	2.0	2.0	72	1	03/21/06	03/21/06	
Ammonia-NH3	EPA 350.3	6C27083	0.14	0.60	ND	1	03/27/06	03/27/06	
Alkalinity as CaCO3	EPA 310.1	6C24119	2.0	2.0	440	1	03/24/06	03/24/06	
Chloride	EPA 300.0	6C21049	3.0	10	220	20	03/21/06	03/21/06	
Nitrate-NO3	EPA 300.0	6C21049	0.35	0.50	8.0	1	03/21/06	03/21/06	
Nitrite-NO2	EPA 300.0	6C21049	6.0	10	ND	20	03/21/06	03/21/06	RL-1
Orthophosphate - PO4	EPA 300.0	6C21049	0.40	0.50	ND	1	03/21/06	03/21/06	
Sulfate	EPA 300.0	6C21049	9.0	10	69	20	03/21/06	03/21/06	
Sulfide	EPA 376.2	6C22079	0.010	0.10	ND	1	03/22/06	03/22/06	
Total Organic Carbon	EPA 415.1	6C24062	0.25	1.0	2.3	1	03/23/06	03/23/06	

### Sample ID: IPC2190-04 (IRZB0095\_WG032106\_0001 - Water)

Reporting Units: mg/l	SM4500-CO2	6C21139	2.0	2.0	55	1	03/21/06	03/21/06	
Carbon Dioxide	SM4500-CO2	6C21139	2.0	2.0	1.1	1	03/21/06	03/21/06	
Ammonia-NH3	EPA 350.3	6C27083	0.14	0.60	640	1	03/24/06	03/24/06	
Alkalinity as CaCO3	EPA 310.1	6C24119	2.0	2.0	120	20	03/21/06	03/21/06	
Chloride	EPA 300.0	6C21049	3.0	10	1.3	1	03/21/06	03/21/06	
Nitrate-NO3	EPA 300.0	6C21049	0.35	0.50	ND	1	03/21/06	03/21/06	
Nitrite-NO2	EPA 300.0	6C21049	0.30	0.50	ND	1	03/21/06	03/21/06	
Orthophosphate - PO4	EPA 300.0	6C21049	0.40	0.50	ND	1	03/21/06	03/21/06	
Sulfate	EPA 300.0	6C21049	9.0	10	79	20	03/21/06	03/21/06	
Sulfide	EPA 376.2	6C22079	0.010	0.10	0.014	1	03/22/06	03/22/06	J
Total Organic Carbon	EPA 415.1	6C24062	0.25	1.0	4.0	1	03/23/06	03/23/06	

### Sample ID: IPC2190-06 (IRZCMW001\_WG032106\_0001 - Water)

Reporting Units: mg/l	EPA 376.2	6C22079	0.010	0.10	ND	1	03/22/06	03/22/06	
Sulfide	EPA 376.2	6C22079	0.010	0.10	ND	1	03/22/06	03/22/06	

### Sample ID: IPC2190-08 (IRZCMW003B\_WG032106\_0001 - Water)

Reporting Units: mg/l	EPA 376.2	6C22079	0.010	0.10	ND	1	03/22/06	03/22/06	
Sulfide	EPA 376.2	6C22079	0.010	0.10	ND	1	03/22/06	03/22/06	

### Sample ID: IPC2190-09 (WCC\_6S\_WG032106\_0001 - Water)

Reporting Units: mg/l	SM4500-CO2	6C21139	2.0	2.0	140	1	03/21/06	03/21/06	
Carbon Dioxide	SM4500-CO2	6C21139	0.14	0.60	0.52	1	03/21/06	03/21/06	J
Ammonia-NH3	EPA 350.3	6C27083	2.0	2.0	520	1	03/24/06	03/24/06	
Alkalinity as CaCO3	EPA 310.1	6C24119	2.0	2.0	460	20	03/21/06	03/22/06	
Chloride	EPA 300.0	6C21049	3.0	10	ND	1	03/21/06	03/22/06	
Nitrate-NO3	EPA 300.0	6C21049	0.35	0.50	ND	20	03/21/06	03/22/06	RL-1
Nitrite-NO2	EPA 300.0	6C21049	6.0	10	ND	1	03/21/06	03/22/06	
Orthophosphate - PO4	EPA 300.0	6C21049	0.40	0.50	ND	1	03/21/06	03/22/06	
Sulfate	EPA 300.0	6C21049	0.45	0.50	12	1	03/21/06	03/22/06	
Total Organic Carbon	EPA 415.1	6C24062	0.25	1.0	17	1	03/23/06	03/23/06	

**Del Mar Analytical - Irvine**  
 Michele Chamberlin  
 Project Manager

L-20750-L

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LDC #: 14879B6  
SDG #: IPC2190  
Laboratory: Del Mar Analytical

## VALIDATION COMPLETENESS WORKSHEET

Tier 2

Date: 5/16/06

Page: 1 of 1

Reviewer: MM

2nd Reviewer: K

**METHOD:** (Analyte) Alkalinity (EPA Method 310.1), Ammonia-X (EPA Method 350.3), Chloride, Nitrate-X, Nitrite-X, Orthophosphate-X, Sulfate (EPA Method 300.0), Sulfide (EPA Method 376.2), TOC (EPA Method 415.1), Carbon Dioxide (SM4500) (CO<sub>2</sub>)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	<b>Validation Area</b>		<b>Comments</b>
I.	Technical holding times	A	Sampling dates: 3/21/06
IIa.	Initial calibration	A	
IIb.	Calibration verification	A	
III.	Blanks	SW X	
IV	Matrix Spike/Matrix Spike Duplicates	A	EPC-190-3 MS/MSB no chrt
V	Duplicates	X A	Dup EPC-190-3
VI.	Laboratory control samples	A	LCS
VII.	Sample result verification	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X	Field blanks	N	

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:

1	IRZB0095_WG032106_0001	11		21		31	
2	P3	12		22		32	
3		13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes:

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LDC #: 14879Bb  
SDG #: IPC2190

## **VALIDATION FINDINGS WORKSHEET**

### **Sample Specific Analysis Reference**

Page: 1 of 1  
Reviewer: my  
2nd reviewer: R

All circled methods are applicable to each sample.

#### **Comments:**

LDC #: 44B19B6  
SDG #: 2PCLYF0

# VALIDATION FINDINGS WORKSHEET

## Blanks

## METHOD: Inorganic Method

See you

Please see qualifications below for all questions answered "N". Not applicable

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".  
 N NA Were all samples associated with a given method blank?  
 N NA Were any inorganic contaminants detected above the reporting limit in the method blanks? If yes, please see qualifications below.

Conc. units: mg/l

## Associated Samples:

41 (358-)

M11 (75X)

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:  
All contaminants within five times the methoc blank concentration were qualified as not detected, "n.d.".

BLANKS. 8

## **Laboratory Data Consultants, Inc. Data Validation Report**

**Project/Site Name:** Boeing Realty Corp, Former C-6 Facility, Torrance CA  
**Collection Date:** March 22, 2006  
**LDC Report Date:** May 15, 2006  
**Matrix:** Water  
**Parameters:** Wet Chemistry  
**Validation Level:** Tier 3  
**Laboratory:** Del Mar Analytical  
**Sample Delivery Group (SDG):** IPC2325

### **Sample Identification**

CMW0002\_WG032206\_0001

## **Introduction**

This data review covers one water sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 300.0 for Chloride, Nitrate, Nitrite, Orthophosphate, and Sulfate, EPA Method 310.1 for Alkalinity, EPA Method 350.3 for Ammonia, EPA Method 376.2 for Sulfide, EPA Method 415.1 for Total Organic Carbon, and Standard Method 4500-CO<sub>2</sub> for Carbon Dioxide.

The review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (October 2004) as there are no current guidelines for the methods stated above.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified a P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blank results are summarized in Section III.

Field duplicates are summarized in Section IX.

The following are definitions of the data qualifiers:

- U     Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
  - J     Indicates an estimated value.
  - R     Quality control indicates the data is not usable.
  - N     Presumptive evidence of presence of the constituent.
  - UJ    Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
  - A     Indicates the finding is based upon technical validation criteria.
  - P     Indicates the finding is related to a protocol/contractual deviation.
- None    Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

All criteria for the initial calibration of each method were met.

### **b. Calibration Verification**

Calibration verification frequency and analysis criteria were met for each method when applicable.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No contaminant concentrations were found in the method blanks.

## **IV. Matrix Spike/Matrix Spike Duplicates**

Matrix spike (MS) and matrix spike duplicate (MSD) analyses were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **V. Duplicates**

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable. Relative percent differences (RPD) were within QC limits.

## **VI. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits.

## **VII. Sample Result Verification**

All sample result verifications were acceptable.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

## **X. Field Blanks**

No field blanks were identified in this SDG.

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Wet Chemistry - Data Qualification Summary - SDG IPC2325**

No Sample Data Qualified in this SDG

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG IPC2325**

No Sample Data Qualified in this SDG



**Del Mar Analytical**

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TAIT Environmental/Boeing  
 701 N. Parkcenter Drive  
 Santa Ana, CA 92705  
 Attention: Mehmet Pehlivan

Project ID: Boeing C-6 Torrance  
 EM2303D/PHASE 01  
 Report Number: IPC2325

Sampled: 03/22/06  
 Received: 03/22/06

## INORGANICS

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
<b>Sample ID: IPC2325-04 (IRZCMW002_WG032206_0001 - Water)</b>									
Reporting Units: mg/l									
Carbon Dioxide	SM4500-CO2	6C23143	2.0	2.0	180	1	03/23/06	03/23/06	
Ammonia-NH3	EPA 350.3	6C29112	0.14	0.60	0.14	1	03/29/06	03/29/06	J
Alkalinity as CaCO3	EPA 310.1	6C27104	2.0	2.0	1000	1	03/27/06	03/27/06	
Chloride	EPA 300.0	6C22049	7.5	25	190	50	03/22/06	03/22/06	
Nitrate-NO3	EPA 300.0	6C22049	0.35	0.50	1.2	1	03/22/06	03/22/06	
Nitrite-NO2	EPA 300.0	6C22049	15	25	35	50	03/22/06	03/22/06	
Orthophosphate - PO4	EPA 300.0	6C22049	0.40	0.50	ND	1	03/22/06	03/22/06	
Sulfate	EPA 300.0	6C22049	0.45	0.50	11	1	03/22/06	03/22/06	
Sulfide	EPA 376.2	6C23091	0.010	0.10	0.035	1	03/23/06	03/23/06	J
Total Organic Carbon	EPA 415.1	6C31057	12	50	510	50	03/30/06	03/30/06	
<b>Sample ID: IPC2325-06 (IRZMW004_WG032206_0001 - Water)</b>									
Reporting Units: mg/l									
Sulfide	EPA 376.2	6C23091	0.010	0.10	ND	1	03/23/06	03/23/06	
<b>Sample ID: IPC2325-07 (IRZMW004_WG032206_0002 - Water)</b>									
Reporting Units: mg/l									
Sulfide	EPA 376.2	6C23091	0.010	0.10	ND	1	03/23/06	03/23/06	
<b>Sample ID: IPC2325-08 (IRZMW002A_WG032206_0001 - Water)</b>									
Reporting Units: mg/l									
Sulfide	EPA 376.2	6C23091	0.010	0.10	ND	1	03/23/06	03/23/06	
<b>Sample ID: IPC2325-09 (CMW0002_WG032206_0001 - Water)</b>									
Reporting Units: mg/l									
Carbon Dioxide	SM4500-CO2	6C23143	2.0	2.0	12	1	03/23/06	03/23/06	
Ammonia-NH3	EPA 350.3	6C29112	0.14	0.60	0.14	1	03/29/06	03/29/06	J
Alkalinity as CaCO3	EPA 310.1	6C27104	2.0	2.0	190	1	03/27/06	03/27/06	
Chloride	EPA 300.0	6C22049	1.5	5.0	120	10	03/22/06	03/22/06	
Nitrate-NO3	EPA 300.0	6C22049	0.35	0.50	ND	1	03/22/06	03/22/06	
Nitrite-NO2	EPA 300.0	6C22049	0.30	0.50	ND	1	03/22/06	03/22/06	
Orthophosphate - PO4	EPA 300.0	6C22049	0.40	0.50	ND	1	03/22/06	03/22/06	
Sulfate	EPA 300.0	6C22049	4.5	5.0	91	10	03/22/06	03/22/06	
Sulfide	EPA 376.2	6C23091	0.010	0.10	0.014	1	03/23/06	03/23/06	J
Total Organic Carbon	EPA 415.1	6C31057	0.25	1.0	15	1	03/30/06	03/30/06	

**Del Mar Analytical - Irvine**  
 Michele Chamberlin  
 Project Manager

125706

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LDC #: 14879C6

SDG #: IPC2325

Laboratory: Del Mar Analytical

## VALIDATION COMPLETENESS WORKSHEET

Tier 3

Date: 5/15/06

Page: 1 of 1

Reviewer: my

2nd Reviewer:

**METHOD: (Analyte)** Alkalinity (EPA Method 310.1), Ammonia-N (EPA Method 350.3), Chloride, Nitrate-N, Nitrite-N, Orthophosphate-N, Sulfate (EPA Method 300.0), Sulfide (EPA Method 376.2), TOC (EPA Method 415.1), Carbon Dioxide (SM4500)  $\text{CO}_2$

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 3/22/06
IIa.	Initial calibration	A	
IIb.	Calibration verification	A	
III.	Blanks	A	
IV	Matrix Spike/Matrix Spike Duplicates	A	3 hrs / 6 hrs / up. was clean +
V	Duplicates	A	
VI.	Laboratory control samples	A	leg
VII.	Sample result verification	A	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X	Field blanks	N	

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:

1	CMW0002_WG032206_0001	11		21		31	
2	MB - P	12		22		32	
3		13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

LDC #: 14879c6  
SDG #: TPC 2325

### VALIDATION FINDINGS CHECKLIST

Page: 1 of 1  
Reviewer: my  
2nd Reviewer: /

Method: Inorganics (EPA Method See copy)

Validation Area	Yes	No	NA	Findings/Comments
<b>I. Technical holding times:</b>				
All technical holding times were met.	✓			
Coolcr tempcrature critcira was met.	✓			
<b>II. Calibration:</b>				
Were all instruments calibrated daily, each set-up time?	✓			
Were the proper number of standards used?	✓			
Were all initial calibration correlation coefficients > 0.995?	✓			
Were all initial and continuing calibration verification %Rs within the 90-110% QC limits?	✓			
Were titrant checks performed as required? (Level IV only)	✓			
Were balance checks performed as required? (Level IV only)			✓	
<b>III. Blanks:</b>				
Was a method blank associated with every sample in this SDG?	✓			
Was there contamination in the method blanks? If yes, please see the Blanks validation completeness worksheet.		✓		
<b>IV. Matrix Spike/Matrix spike duplicates and Duplicates:</b>				
Were a matrix spike (MS) and duplicate (DUP) analyzed for each matrix in this SDG? If no, indicate which matrix does not have an associated MS/MSD or MS/DUP. Soil / Water.	✓			<i>no chart</i>
Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the 75-125 QC limits? If the sample concentration exceeded the spike concentration by a factor of 4 or more, no action was taken.	✓			
Were the MS/MSD or duplicate relative percent differences (RPD) $\leq$ 20% for waters and $\leq$ 35% for soil samples? A control limit of $\leq$ CRDL ( $\leq$ 2X CRDL for soil) was used for samples that were $\leq$ 5X the CRDL, including when only one of the duplicate sample values were $<$ 5X the CRDL.	✓			
<b>V. Laboratory control samples:</b>				
Was an LCS analyzed for this SDG?	✓			
Was an LCS analyzed per extraction batch?	✓			
Were the LCS percent recoveries (%R) and relative percent difference (RPD) within the 80-120% (85-115% for Method 300.0) QC limits?	✓			
<b>VI. Regional Quality Assurance and Quality Control:</b>				
Were performance evaluation (PE) samples performed?			✓	
Were the performance evaluation (PE) samples within the acceptance limits?			✓	

LDC #: 14879cb  
SDG #: TPLC3-N

VALIDATION FINDINGS CHECKLIST

Page: 1 of 1  
Reviewer: my  
2nd Reviewer: K

Validation Area	Yes	No	NA	Findings/Comments
<b>VII. Sample Result Verification</b>				
Were RLs adjusted to reflect all sample dilutions and dry weight factors applicable to level IV validation?	✓			
Were detection limits < RL?	✓			
<b>VIII. Overall assessment of data</b>				
Overall assessment of data was found to be acceptable.	✓			
<b>IX. Field duplicates</b>				
Field duplicate pairs were identified in this SDG.		✓		
Target analytes were detected in the field duplicates.			✓	
<b>X. Field blanks</b>				
Field blanks were identified in this SDG.		✓		
Target analytes were detected in the field blanks.			✓	

LDC #: 14879cb  
SDG #: 1P623V5

**VALIDATION FINDINGS WORKSHEET**  
Sample Specific Analysis Reference

Page: 1 of 1  
Reviewer: WY  
2nd reviewer:   

All circled methods are applicable to each sample.

Sample ID	Parameter
1	pH TDS <u>C</u> F <u>NO<sub>3</sub></u> <u>NO<sub>2</sub></u> <u>SO<sub>4</sub></u> <u>PO<sub>4</sub></u> <u>ALK</u> CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> <u>(C)</u> <u>(S)</u> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____
	pH TDS Cl F NO <sub>3</sub> NO <sub>2</sub> SO <sub>4</sub> PO <sub>4</sub> ALK CN <sup>-</sup> NH <sub>3</sub> TKN TOC CR <sup>δ+</sup> _____

Comments: \_\_\_\_\_

LDC #: 14879 C6  
SDG #: TC238

Validation Findings Worksheet  
Initial and Continuing Calibration Calculation Verification

Page: 1 of 1  
Reviewer: WJ  
2nd Reviewer: RC

Method: Inorganics, Method See con

The correlation coefficient ( $r$ ) for the calibration of SO<sub>4</sub> was recalculated for each type of analysis using the following formula:

$$\%R = \frac{\text{Found} \times 100}{\text{True}}$$

Where,  
Found = concentration of each analyte measured in the analysis of the ICV or CCV solution  
True = concentration of each analyte in the ICV or CCV source

An initial or continuing calibration verification percent recovery (%R) was recalculated for each type of analysis using the following formula:

Type of analysis	Analyte	Standard	Conc. (mg/L)	Area	Recalculated	Reported	Acceptable
Initial calibration	SO <sub>4</sub>	s1	0	0			
Calibration verification		s2	0.4	37086.13	0.999994	0.999994	
		s3	1	78840.36			
		s4	10	882302.78			
		s5	20	1859685.6			
		s6	40	4073330.88			
		s7	60	6589508.82			
Calibration verification	Cl	10	9.8011	9.8	9.8	9.8	
Calibration verification	Na <sub>3</sub>	4	4.08	4.08	4.08	4.08	
Calibration verification	T <sub>OC</sub>	60	9.023	9.023	9.023	9.023	

Comments: Refer to Calibration Verification findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

LDC #: 14879C6  
SDG #: TPU-335

**VALIDATION FINDINGS WORKSHEET**  
**Level IV Recalculation Worksheet**

Page: 1 of 1  
Reviewer: MM  
2nd Reviewer: EC

METHOD: Inorganics, Method See Curve

Percent recoveries (%R) for a laboratory control sample and a matrix spike sample were recalculated using the following formula:

%R =  $\frac{\text{Found}}{\text{True}} \times 100$  Where, Found = concentration of each analyte measured in the analysis of the sample. For the matrix spike calculation,  
Found = SSR (spiked sample result) - SR (sample result).  
True = concentration of each analyte in the source.

A sample and duplicate relative percent difference (RPD) was recalculated using the following formula:

$$\text{RPD} = \frac{|S-D|}{(S+D)/2} \times 100 \quad \text{Where, } S = \text{Original sample concentration}$$
$$D = \text{Duplicate sample concentration}$$

Sample ID	Type of Analysis	Element	Found / S (units)	True / D (units)	Recalculated		Reported %R / RPD	Acceptable (Y/N)
					%R / RPD	Reported		
Lc3	Laboratory control sample	Co <sub>2</sub>	4.89	5.0	98	98	Y	
TPU-1	Matrix spike sample	S	0.467	0.580	81	80	J	
TPU-4	Duplicate sample	Co <sub>2</sub>	180	180	0	0	C	

Comments: Refer to appropriate worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

LDC #: 14879cb  
SDG #: IPC2325

## **VALIDATION FINDINGS WORKSHEET**

### Sample Calculation Verification

Page: 1 of 1  
Reviewer: ky  
2nd reviewer: N

**METHOD:** Inorganics, Method See cover

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

- N N/A Have results been reported and calculated correctly?  
 N N/A Are results within the calibrated range of the instruments?  
 N N/A Are all detection limits below the CRQL?

Compound (analyte) results for \_\_\_\_\_ reported with a positive detect were recalculated and verified using the following equation:

**Concentration =**

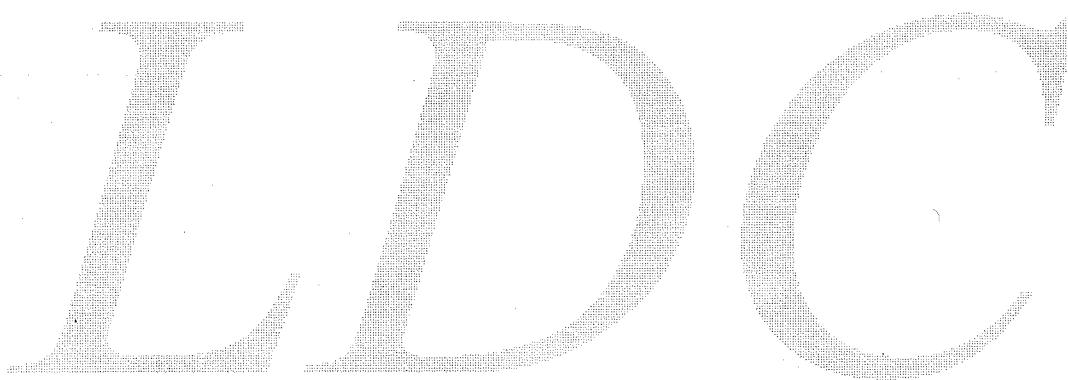
### Recalculation:

$$\text{Alkalinity} = \frac{\text{NaOH} \times \sqrt{\text{Lactate}} \times 5000}{\text{Sample Volume}} \quad \text{Alkalinity} = \frac{4.70 \times 0.02 \times 5000}{25.0} = 188 \text{ mg/L}$$

Note: \_\_\_\_\_

**Boeing Realty Corp., Former C-6 Facility, Torrance CA**  
**Data Validation Reports**  
**LDC# 14879**

**Fixed Gases**



## **Laboratory Data Consultants, Inc.**

### **Data Validation Report**

**Project/Site Name:** Boeing Realty Corp, Former C-6 Facility, Torrance CA

**Collection Date:** March 20, 2006

**LDC Report Date:** May 15, 2006

**Matrix:** Water

**Parameters:** Fixed Gases

**Validation Level:** Tier 1

**Laboratory:** Del Mar Analytical\Air Technology Laboratory, Inc.

**Sample Delivery Group (SDG):** IPC2080/A6032106

#### **Sample Identification**

MWC011\_WG032006\_0001

## **Introduction**

This data review covers one water sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per Method RSK-175 for Fixed Gases.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review (October 1999) as there are no current guidelines for the method stated above.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified a P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blank results are summarized in Section III.

Field duplicates are summarized in Section IX.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- U      Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
  - J      Indicates an estimated value.
  - R      Quality control indicates the data is not usable.
  - N      Presumptive evidence of presence of the constituent.
  - UJ     Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
  - A      Indicates the finding is based upon technical validation criteria.
  - P      Indicates the finding is related to a protocol/contractual deviation.
- None    Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures.

The cooler temperatures for sample MWC011\_WG032006\_0001 was reported at 10°C upon receipt by the laboratory.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration data were not reviewed for Tier 1.

### **b. Calibration Verification**

Calibration verification data were not reviewed for Tier 1.

## **III. Blanks**

Method blanks were performed at the required frequency. No fixed gases contaminants were found in the method blanks.

## **IV. Accuracy and Precision Data**

### **a. Surrogate Recovery**

Surrogates were not required by the method.

### **b. Matrix Spike/Matrix Spike Duplicates**

Matrix spike (MS) and matrix spike duplicate (MSD) analyses were not required by the method.

### **c. Laboratory Control Samples**

Laboratory control samples were analyzed at the required frequency. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **V. Target Compound Identification**

Raw data were not reviewed for this SDG.

## **VI. Compound Quantitation and CRQLs**

Raw data were not reviewed for this SDG.

## **VII. System Performance**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

## **X. Field Blanks**

No field blanks were identified in this SDG.

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Fixed Gases - Data Qualification Summary - SDG IPC2080/A6032106**

No Sample Data Qualified in this SDG

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Fixed Gases - Laboratory Blank Data Qualification Summary - SDG**  
**IPC2080/A6032106**

No Sample Data Qualified in this SDG

Client: Del Mar Analytical, Irvine  
Attn: Michele Chamberlin

Page 2 of 3  
A6032106

Client's Project: IPC2080  
Date Received: 3/21/2006  
Matrix: Water  
Units: ug/L

Dissolved Gases by EPA Procedure RSKSOP-175

Lab No.:	A6032106-01	A6032106-02			
Client Sample I.D.:	IPC2080-07	IPC2080-11			
Date Sampled:	3/20/2006	3/20/2006			
Date Analyzed:	3/27/2006	3/27/2006			
Analyst Initials:	DT	DT			
Data File:	27mar011	27mar012			
QC Batch:	060327GC8A1	060327GC8A1			
Dilution Factor:	1.0	1.0			
ANALYTE	PQL	RL	Results	RL	Results
Methane	1.0	1.0	88	1.0	ND
Ethane	2.0	2.0	ND	2.0	ND
Ethylene	3.0	3.0	ND	3.0	ND
Nitrogen	1500	1500	110,000	1500	97,000

PQL = Practical Quantitation Limit

ND = Not Detected (Below RL)

RL = PQL X Dilution Factor

Reviewed/Approved By:

Date: 3-31-06

Mark J. Johnson  
Operations Manager

The cover letter is an integral part of this analytical report.



AirTECHNOLOGY Laboratories, Inc.

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LDC #: 14879A51

**VALIDATION COMPLETENESS WORKSHEET**

Date: 5/12/06

SDG #: IPC2080/A6032106

Tier 1

Page: 1 of 1

Laboratory: Del Mar Analytical/Air Technology Laboratory, Inc.

Reviewer: P

2nd Reviewer: u

**METHOD: GC Fixed Gases (Method RSK-175)**

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	<b>Validation Area</b>		<b>Comments</b>
I.	Technical holding times	S/N	Sampling dates: 3/20/06
IIa.	Initial calibration	N	
IIb.	Calibration verification	N	
III.	Blanks	Δ	
IVa.	Surrogate recovery	A	not required
IVb.	Matrix spike/Matrix spike duplicates	N	client specified
IVc.	Laboratory control samples	A	LCS ID
V.	Target compound identification	N	
VI.	Compound Quantitation and CRQLs	N	
VII.	System Performance	N	
VIII.	Overall assessment of data	Δ	
IX.	Field duplicates	N	
X.	Field blanks	N	

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet

ND = No compounds detected  
 R = Rinsate  
 FB = Field blank

D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples: water

+	MWC011_WG032006_0001	11	Method Bk 3/21/06	21		31	
2		12		22		32	
3		13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes:

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LDC #: 14879A51  
SDG #: PC2080/A6032104

# VALIDATION FINDINGS WORKSHEET

## Technical Holding Times

All circled dates have exceeded the technical holding times.  
Y N N/A Were all cooler temperatures within validation criteria?

Page: \_\_\_\_\_  
Reviewer: \_\_\_\_\_  
2nd Reviewer: \_\_\_\_\_

Reviewer: \_\_\_\_\_

2nd Reviewer:

卷之三

## METHOD: GC HPLC

TECHNICAL HOLDING TIME CRITERIA

**VOLATILES:** Water unpreserved:

Water preserved:  
Soils:

SYNTHETIC POLYMERS

3. Water:  
Soil:

**A** Aromatic within 7 days, non-aromatic within 14 days of sample collection.

Both within 14 days of sample collection.

Extracted within 7 days, analyzed within 40 days.  
Extracted within 14 days, analyzed within 40 days

## **Laboratory Data Consultants, Inc. Data Validation Report**

**Project/Site Name:** Boeing Realty Corp, Former C-6 Facility, Torrance CA  
**Collection Date:** March 21, 2006  
**LDC Report Date:** May 15, 2006  
**Matrix:** Water  
**Parameters:** Fixed Gases  
**Validation Level:** Tier 2  
**Laboratory:** Del Mar Analytical\Air Technology Laboratory, Inc.  
**Sample Delivery Group (SDG):** IPC2190/A6032203

### **Sample Identification**

IRZB0095\_WG032106\_0001

## **Introduction**

This data review covers one water sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per Method RSK-175 for Fixed Gases.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review (October 1999) as there are no current guidelines for the method stated above.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified a P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blank results are summarized in Section III.

Field duplicates are summarized in Section IX.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- U     Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
  - J     Indicates an estimated value.
  - R     Quality control indicates the data is not usable.
  - N     Presumptive evidence of presence of the constituent.
  - UJ    Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
  - A     Indicates the finding is based upon technical validation criteria.
  - P     Indicates the finding is related to a protocol/contractual deviation.
- None    Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## I. Technical Holding Times

All technical holding time requirements were met.

The chain-of-custodices were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria with the following exceptions:

Sample	Compound	Finding	Criteria	Flag	A or P
IRZB0095_WG032106_0001	All TCL compounds	Cooler temperature was reported at 11°C upon receipt by the laboratory.	Cooler temperature must be $4\pm2^{\circ}\text{C}$ .	J (all detects) UJ (all non-detects)	A

## II. Calibration

### a. Initial Calibration

Initial calibration of compounds was performed as required by the method.

A curve fit, based on the initial calibration, was established for quantitation. The coefficient of determination ( $r^2$ ) was greater than or equal to 0.990 .

### b. Calibration Verification

Calibration verification was performed at required frequencies. The percent differences (%D) of amounts in continuing standard mixtures were within the 25.0% QC limits.

## III. Blanks

Method blanks were performed at the required frequency. No fixed gases contaminants were found in the method blanks.

## IV. Accuracy and Precision Data

### a. Surrogate Recovery

Surrogates were not required by the method.

### b. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) analyses were not required by the method.

### c. Laboratory Control Samples

Laboratory control samples were analyzed at the required frequency. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **V. Target Compound Identification**

Raw data were not reviewed for this SDG.

## **VI. Compound Quantitation and CRQLs**

Raw data were not reviewed for this SDG.

## **VII. System Performance**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

## **X. Field Blanks**

No field blanks were identified in this SDG.

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Fixed Gases - Data Qualification Summary - SDG IPC2190/A6032203**

SDG	Sample	Compound	Flag	A or P	Reason
IPC2190/ A6032203	IRZB0095_WG032106_0001	All TCL compounds	J (all detects) UJ (all non-detects)	A	Cooler temperature

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Fixed Gases - Laboratory Blank Data Qualification Summary - SDG**  
**IPC2190/A6032203**

No Sample Data Qualified in this SDG

Client: Del Mar Analytical, Irvine  
Attn: Michele Chamberlin

Page 2 of 3  
A6032203

Client's Project: IPC2190  
Date Received: 3/22/2006  
Matrix: Water  
Units: ug/L

Dissolved Gases by EPA Procedure RSKSOP-175

Lab No.:	A6032203-01	A6032203-02	A6032203-03	A6032203-04	
Client Sample I.D.:	IPC2190-03	IPC2190-04	IPC2190-09	IPC2190-11	
Date Sampled:	3/21/2006	3/21/2006	3/21/2006	3/21/2006	
Date Analyzed:	3/27/2006	3/27/2006	3/27/2006	3/27/2006	
Analyst Initials:	DT	DT	DT	DT	
Data File:	27mar013	27mar014	27mar015	27mar016	
QC Batch:	060327GC8A1	060327GC8A1	060327GC8A1	060327GC8A1	
Dilution Factor:	1.0	1.0	1.0	1.0	
ANALYTE	PQL	RL	Results	RL	Results
Methane	1.0	1.0	1,800	1.0	12,000
Ethane	2.0	2.0	ND	2.0	ND
Ethylene	3.0	3.0	ND	3.0	4.1
Nitrogen	1500	1500	97,000	1500	82,000
				1500	94,000
				1500	110,000

PQL = Practical Quantitation Limit

ND = Not Detected (Below RL)

RL = PQL X Dilution Factor

Reviewed/Approved By:

Mark J. Johnson  
Operations Manager

Date: 3-31-06

The cover letter is an integral part of this analytical report.



AirTECHNOLOGY Laboratories, Inc.

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LDC #: 14879B51**VALIDATION COMPLETENESS WORKSHEET**Date: 5/2/06SDG #: IPC2190/A6032203

Tier 2

Page: 1 of 1Laboratory: Del Mar Analytical/Air Technology Laboratory, Inc.Reviewer: J2nd Reviewer: R**METHOD: GC Fixed Gases (Method RSK-175)**

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	<b>Validation Area</b>		<b>Comments</b>
I.	Technical holding times	S/N	Sampling dates: <u>3/21/06</u>
IIa.	Initial calibration	A	<u>% PSP &lt; 25</u> ✓ <u>± 0.990</u>
IIb.	Calibration verification	A	<u>% D &lt; 25</u>
III.	Blanks	A	
IVa.	Surrogate recovery	N	not required
IVb.	Matrix spike/Matrix spike duplicates	N	chint specified
IVc.	Laboratory control samples	A	LCS ID
V.	Target compound identification	N	
VI.	Compound Quantitation and CRQLs	N	
VII.	System Performance	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X.	Field blanks	N	

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet

ND = No compounds detected  
 R = Rinsate  
 FB = Field blank

D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples:

water

1	IRZB0095_WG032106_0001	11	Method Blank 3/21/06	21		31	
2		12		22		32	
3		13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes:

LDC #: 14879B5 /  
SDG #: 1PC2190 / A6032203

## **VALIDATION FINDINGS WORKSHEET**

Technical Holding Times

## ATION FINDINGS WORKS Technical Holding Times

All circled dates have exceeded the technical holding times.  
Y N NA Were all cooler temperatures within validation crit.

## METHOD: ✓GC HPLC

TECHNICAL HOLDING TIME CRITERIA

**VOLATILES:** Water unpreserved:

Water preserved:

## Soils:

Water:  
Soil:

-S.  
Water:  
soil:

Aromatic: within 7 days non-aromatic within 14 days of sample collection

Both within 14 days of sample collection

Extracted within 7 days, analyzed within 40 days.

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** Boeing Realty Corp, Former C-6 Facility, Torrance CA  
**Collection Date:** March 22, 2006  
**LDC Report Date:** May 15, 2006  
**Matrix:** Water  
**Parameters:** Fixed Gases  
**Validation Level:** Tier 3  
**Laboratory:** Del Mar Analytical\Air Technology Laboratory, Inc.  
**Sample Delivery Group (SDG):** IPC2325/A6032303

**Sample Identification**

CMW0002\_WG032206\_0001

## **Introduction**

This data review covers one water sample listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per Method RSK-175 for Fixed Gases.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review (October 1999) as there are no current guidelines for the method stated above.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified a P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blank results are summarized in Section III.

Field duplicates are summarized in Section IX.

The following are definitions of the data qualifiers:

- U     Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- J     Indicates an estimated value.
- R     Quality control indicates the data is not usable.
- N     Presumptive evidence of presence of the constituent.
- UJ    Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A     Indicates the finding is based upon technical validation criteria.
- P     Indicates the finding is related to a protocol/contractual deviation.
- None   Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures.

The cooler temperature for sample CMW0002\_WG032206\_000 was reported at 10°C upon receipt by the laboratory.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration of compounds was performed as required by the method.

A curve fit, based on the initial calibration, was established for quantitation. The coefficient of determination ( $r^2$ ) was greater than or equal to 0.990 .

### **b. Calibration Verification**

Calibration verification was performed at required frequencies. The percent differences (%D) of amounts in continuing standard mixtures were within the 25.0% QC limits.

## **III. Blanks**

Method blanks were performed at the required frequency. No fixed gases contaminants were found in the method blanks.

## **IV. Accuracy and Precision Data**

### **a. Surrogate Recovery**

Surrogates were not required by the method.

### **b. Matrix Spike/Matrix Spike Duplicates**

Matrix spike (MS) and matrix spike duplicate (MSD) analyses were not required by the method.

### **c. Laboratory Control Samples**

Laboratory control samples were analyzed at the required frequency. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

## **V. Target Compound Identification**

All target compound identifications were within validation criteria.

## **VI. Compound Quantitation and CRQLs**

All compound quantitation and CRQLs were within validation criteria.

## **VII. System Performance**

The system performance was acceptable.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report if data has been qualified.

## **IX. Field Duplicates**

No field duplicates were identified in this SDG.

## **X. Field Blanks**

No field blanks were identified in this SDG.

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Fixed Gases - Data Qualification Summary - SDG IPC2325/A6032303**

No Sample Data Qualified in this SDG

**Boeing Realty Corp, Former C-6 Facility, Torrance CA**  
**Fixed Gases - Laboratory Blank Data Qualification Summary - SDG**  
**IPC2325/A6032303**

No Sample Data Qualified in this SDG

Client: Del Mar Analytical, Irvine  
Attn: Michele Chamberlin

Page 2 of 3  
A6032303

Client's Project: IPC2325  
Date Received: 3/23/2006  
Matrix: Water  
Units: ug/L

Dissolved Gases by EPA Procedure RSKSOP-175

Lab No.:	A6032303-01	A6032303-02			
Client Sample I.D.:	IPC2325-04	IPC2325-09			
Date Sampled:	3/22/2006	3/22/2006			
Date Analyzed:	3/27/2006	3/27/2006			
Analyst Initials:	DT	DT			
Data File:	27mar017	27mar018			
QC Batch:	060327GC8A1	060327GC8A1			
Dilution Factor:	1.0	1.0			
ANALYTE	PQL	RL	Results	RL	Results
Methane	1.0	1.0	12,000	1.0	1.6
Ethane	2.0	2.0	ND	2.0	ND
Ethylene	3.0	3.0	ND	3.0	ND
Nitrogen	1500	1500	52,000	1500	110,000

PQL = Practical Quantitation Limit

ND = Not Detected (Below RL)

RL = PQL X Dilution Factor

Reviewed/Approved By:

Mark J. Johnson  
Operations Manager

Date: 3-31-06

The cover letter is an integral part of this analytical report.



AirTECHNOLOGY Laboratories, Inc.

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LDC #: 14879C51

**VALIDATION COMPLETENESS WORKSHEET**

Date: 5/15/06

SDG #: IPC2325/A6032303

Tier 3

Page: 1 of

Laboratory: Del Mar Analytical/Air Technology Laboratory, Inc.

Reviewer: J

2nd Reviewer: J

**METHOD: GC Fixed Gases (Method RSK-175)**

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	<b>Validation Area</b>		<b>Comments</b>
I.	Technical holding times	SW	Sampling dates: 3/22/06
IIa.	Initial calibration	A	$r^2 = 0.990$
IIb.	Calibration verification	A	$cev \leq 25\%$
III.	Blanks	A	
IVa.	Surrogate recovery	N	not required
IVb.	Matrix spike/Matrix spike duplicates	N	client specified
IVc.	Laboratory control samples	A	lcs ID
V.	Target compound identification	A	
VI.	Compound Quantitation and CRQLs	A	
VII.	System Performance	A	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	N	
X.	Field blanks	N	

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet

ND = No compounds detected  
 R = Rinsate  
 FB = Field blank

D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples:

water

1	CMW0002_WG032206_0001	11	Method Blank 3/27/06	21		31	
2		12		22		32	
3		13		23		33	
4		14		24		34	
5		15		25		35	
6		16		26		36	
7		17		27		37	
8		18		28		38	
9		19		29		39	
10		20		30		40	

Notes:

LDC #: 1487451  
SDG #: 1PC2325/A 6032303

### VALIDATION FINDINGS CHECKLIST

Page: 6 of 2  
Reviewer:       
2nd Reviewer:     

Method: GC HPLC

Validation Area	Yes	No	NA	Findings/Comments
<b>I. Technical holding times</b>				
All technical holding times were met.	✓	1		
Cooler temperature criteria was met.		✓		
<b>II. Initial calibration</b>				
Did the laboratory perform a 5 point calibration prior to sample analysis?	✓			
Was a linear fit used for evaluation? If yes, were all percent relative standard deviations (%RSD) < 20%?			✓	
Was a curve fit used for evaluation? If Yes, what was the acceptance criteria used?	✓			
Did the initial calibration meet the curve fit acceptance criteria?	✓			
Were the RT windows properly established?	✓			
<b>IV. Continuing calibration</b>				
What type of continuing calibration calculation was performed? ___ %D or ___ %R	✓			
Was a continuing calibration analyzed daily?	✓			
Were all percent differences (%D) < 15% or percent recoveries 85-115%?	✓			
Were all the retention times within the acceptance windows?	✓			
<b>V. Blanks</b>				
Was a method blank associated with every sample in this SDG?	✓			
Was a method blank analyzed for each matrix and concentration?	✓			
Was there contamination in the method blanks? If yes, please see the Blanks validation completeness worksheet.		✓		
<b>VI. Surrogate spikes</b>				
Were all surrogate %R within the QC limits?			✓	
If the percent recovery (%R) of one or more surrogates was outside QC limits, was a reanalysis performed to confirm %R?			✓	
If any %R was less than 10 percent, was a reanalysis performed to confirm %R?			✓	
<b>VII. Matrix spike/Matrix spike duplicates</b>				
Were a matrix spike (MS) and matrix spike duplicate (MSD) analyzed for each matrix in this SDG? If no, indicate which matrix does not have an associated MS/MSD. Soil / Water.			✓	
Was a MS/MSD analyzed every 20 samples of each matrix?			✓	
Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the QC limits?			✓	
<b>VIII. Laboratory control samples</b>				
Was an LCS analyzed for this SDG?	✓			
Was an LCS analyzed per extraction batch?	✓			

LDC #: 14879CS1  
SDG #: IPC 2325/A 6032303

### VALIDATION FINDINGS CHECKLIST

Page: 7 of 2  
Reviewer: J  
2nd Reviewer: R

Validation Area	Yes	No	NA	Findings/Comments
Were the LCS percent recoveries (%R) and relative percent difference (RPD) within the QC limits?	/			
<b>X. Regional Quality Assurance and Quality Control</b>				
Were performance evaluation (PE) samples performed?		/		
Were the performance evaluation (PE) samples within the acceptance limits?			/	
<b>XI. Target compound identification</b>				
Were the retention times of reported detects within the RT windows?	/			
<b>XII. Compound quantitation/CRQLs</b>				
Were compound quantitation and CRQLs adjusted to reflect all sample dilutions and dry weight factors applicable to level IV validation?	/			
<b>XIII. System performance</b>				
System performance was found to be acceptable.	/			
<b>XIV. Overall assessment of data</b>				
Overall assessment of data was found to be acceptable.	/			
<b>XV. Field duplicates</b>				
Were field duplicate pairs identified in this SDG?		/		
Were target compounds detected in the field duplicates?			/	
<b>XVI. Field blanks</b>				
Were field blanks identified in this SDG?		/		
Were target compounds detected in the field blanks?			/	

LDC #: 14879c5  
SDG #: 1PC237S/A6032303

# VALIDATION FINDINGS WORKSHEET

## Technical Holding Times

All circled dates have exceeded the technical holding times.

All circled dates have exceeded the technical holding times.

Page: 1 of 1  
Reviewer: J  
2nd Reviewer: J

## METHOD: GC HPLC

TECHNICAL HIGHLIGHTING TIME CRITERIA

**CHINAS** **WALL** **PAINT**  
**WOLTLIES:** Water unpreserved;  
Water preserved:

Both Within 14 days of sample collection.  
Both within 14 days of sample collection.

**EXTRACTABLES:** Water: Soil:

Extracted within 7 days, analyzed within 40 days.  
Extracted within 14 days, analyzed within 40 days.

non-aromatic within 14 days of sample collection.

Multiple collection

Both within 14 days of sample collection.

Both within 14 days of sample collection.

Extracted within 7 days, analyzed within 40 days.  
Extracted within 14 days, analyzed within 40 days.

LDC# 14879c51  
SDG# 1P02325 /A-6032303

**VALIDATION FINDINGS WORKSHEET**  
Initial Calibration Calculation Verification

METHOD: \_\_\_\_\_

8082

Page: 1 of 2  
Reviewer: J  
2nd Reviewer: A

Calibration Date	Column/ Detector	Compound	Standard	X	Y
04/20/05	FID	methane	Point 1	10	13683
			Point 2	100	121591
			Point 3	1000	1260392
			Point 4	5000	6514911
			Point 5	10000	13405678
			Point 6		
			Point 7		
			Point 7		
			Point 8		

Regression Output:	Recalculated Result	Result Reported by the Laboratory
Constant	0	0.00E+00
Std Err of Y Est	91688.222524795768	
R Squared	0.99975	0.999811
No. of Observations	5	
Degrees of Freedom	4	
X Coefficient(s)	1332.46480187	
Std Err of Coef.	8.16790871554	1.3325E+03

LDC# 14879es  
SDG# 110c 2325/A 6032303

**VALIDATION FINDINGS WORKSHEET**  
**Initial Calibration Calculation Verification**

METHOD:       

8082

Page: 2 of 2  
Reviewer: B  
2nd Reviewer:   

Calibration Date	Column/ Detector	Compound	Standard	X	Y
03/13/06	TCD	nitrogen	Point 1	1.56	63916
			Point 2	7.81	290198
			Point 3	39.099998	1393537
			Point 4	100	3442888
			Point 5		
			Point 6		
			Point 7		
			Point 7		
			Point 8		

Recalculated Result		Result Reported by the Laboratory
Regression Output:		
Constant	0	0.00E+00
Std Err of Y Est	28547.659624440576	
R Squared	0.99966	0.999818
No. of Observations	4	
Degrees of Freedom	3	
X Coefficient(s)	34604.3788816	3.4604E+04
Std Err of Coef.	265.147072585	

LDC #: 14879051  
SDG #: 11PC 2325/A 6032 303

**VALIDATION FINDINGS WORKSHEET**  
**Continuing Calibration Results Verification**

Page: 6 of 7  
Reviewer: J  
2nd Reviewer: J

**METHOD: GC** **HPLC**

The percent difference (%D) of the initial calibration average Calibration Factors (CF) and the continuing calibration CF were recalculated for the compounds identified below using the following calculation:

% Difference =  $100 * (\text{ave. CF} - \text{CF})/\text{ave. CF}$   
CF = A/C  
A = Area of compound  
C = Concentration of compound

#	Standard ID	Calibration Date	Compound	Average CF(cal)/CCV Conc.	CF/Conc. CCV	Reported	Recalculated	Reported	Recalculated
						%D	%D	%D	%D
1	nitrobenzene	3/27/06 1:52 AM	nitrobenzene	5000	4467.8	4467.8	10.6	10.6	10.6
2	nitrobenzene	3/27/06 8:21 AM	nitrobenzene	39.10	39.115	39.115	0	0	0
3									
4									

Comments: Refer to Continuing Calibration findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

LDC #: 144-79CS  
SDG #: 1PC-2-3 25/AU03-2303

# VALIDATION FINDINGS WORKSHEET

## Laboratory Control Sample Duplicate Results Verification

Page: 6 of 7  
Reviewer: J  
2nd Reviewer

METHOD: GC HPLC

The percent recoveries (%R) and relative percent differences (RPD) of the laboratory control sample and laboratory control sample duplicate were recalculated for the compounds identified below using the following calculation:

%Recovery = 100 \* (SSC - SC)/SA

Where  
SSC = Spiked concentration  
SA = Spike added

RPD = {{(SSCLCS - SSCLCSD)} \* 2} / (SSCLCS + SSCLCSD)) \* 100

LCS/LCSD samples: 2/2 106

SC = Sample concentration

LCS = Laboratory Control Sample percent recovery

LCSD = Laboratory Control Sample duplicate percent recovery

Compound	Spike Added			Sample Conc.			Spike Sample Concentration			Percent Recovery		Percent Recovery		LCS/LCSD	
	LCS	LCSD	—	LCS	LCSD	—	LCS	LCSD	—	Reported	Recalc.	Reported	Recalc.	Reported	RPD
Gasoline (8015)															
Diesel (8015)															
Benzene (8021B)															
Methane (RSK-175)	6900.0	6900.0	0	6761.3	6516.7	—	97	98	94	94	94	3.7	3.7		
2,4-D (8151)															
Dinitrobenzene (8151)															
Naphthalene (8310)															
Anthracene (8310)															
HMX (8330)															
2,4,6-Trinitrotoluene (8330)															

Comments: Refer to Laboratory Control Sample/Laboratory Control Sample Duplicate findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

LDC #: 1481905  
SDG #: 17C 2325 / A603-2303

**VALIDATION FINDINGS WORKSHEET**  
**Sample Calculation Verification**

Page: 1 of 1  
Reviewer: J  
2nd Reviewer: A

METHOD: GC HPLC

(Y) N/A  
(Y) N/A

Were all reported results recalculated and verified for all level IV samples?  
Were all recalculated results for detected target compounds within 10% of the reported results?

$$\text{Concentration} = \frac{(A)(F_v)(D_f)}{(RF)(V_s \text{ or } W_s)(\%S/100)}$$

A= Area or height of the compound to be measured  
Fv= Final Volume of extract  
Df= Dilution Factor  
RF= Average response factor of the compound  
In the initial calibration  
Vs= Initial volume of the sample  
Ws= Initial weight of the sample  
%S= Percent Solid

Example:

Sample ID. #1

Compound Name Methane

$\frac{1.6}{1.4} = 1.52 \times 100 \approx 5\%$

$$\text{Concentration} = \frac{(0.00037 + 0.0014667)}{1000}$$

$$= 1.52 \text{ ng/L}$$

#	Sample ID	Compound	Reported Concentrations	Recalculated Results Concentrations	Qualifications
#1	Methane	$\frac{1.3325 \times 10^3}{22900} = 1.3325 \times 10^{-3}$ (X)			
		$X = 17.2 \text{ ppm}$			
	in water	$\frac{17.2}{10^6} \times 55.51 \times 16.04 \times \frac{1000}{41300} = 0.00037$			
	in H2O	$17.2 \times 55.06 \times 16.04 \times \frac{1000}{1000} \times \frac{1}{40} = 0.0014667$			

Comments: \_\_\_\_\_